Assigment 5

Upload your code (.ipynb) on Learn dropbox and submit pdfs of the code and to Crowdmark.	

Classification

```
In [880]:
          # Download the LIBSVM package from here: https://www.csie.ntu.edu.tw/
          ~cilin/libsvm/#download
          # If your download is successfull you should have the folder with nam
          e: libsvm-3.24.
          # We will use this package to load datasets.
          # Enter the downloaded folder libsvm-3.24 through your terminal.
          # Run make command to compile the package.
          # Load this auxiliary package.
          import sys
          # add here your path to the folder libsvm-3.24/python
          path = "/home/tempo/Desktop/Fall 2019/CS 794/a5/libsvm-3.24/python"
          # Add the path to the Python paths so Python can find the module.
          sys.path.append(path)
          # Load the LIBSVM module.
          from symutil import *
          # Add here your path to the folder libsvm-3.24
          path = "/home/tempo/Desktop/Fall 2019/CS 794/a5/libsvm-3.24"
          # Test that it works. This will load the data "heart scale"
          # and it will store the labels in "b" and the data matrix in "A".
          b, A = svm read problem(path + '/heart scale')
          print('Loaded data: Heart Scale')
          # Use "svm read problem" function to load data for your assignment.
          # Note that matrix "A" stores the data in a sparse format.
          # In particular matrix "A" is a list of dictionaries.
          # The length of the list gives you the number of samples.
          # Each entry in the list is a dictionary. The keys of the dictionary
           are the non-zero features.
          # The values of the dictionary for each key is a list which gives you
          the feature value.
```

Loaded data: Heart Scale

```
In [881]: import matplotlib.pyplot as plt

# Numpy is useful for handling arrays and matrices.
import numpy as np
from scipy.sparse import coo_matrix
import time
from random import randrange as rnd
import sklearn
import sklearn.metrics
import random
import pandas as pd
```

```
path news20 = r'/home/tempo/Desktop/Fall 2019/CS 794/a5/news20.binar
In [882]:
           b news20, A news20 = svm read problem(path news20) \#(19996, 1355191)
           def splitData(A, b):
               n = A.shape[0]
               test split = 15996
               valid split = 2000
               train split = 2000
               ATrain = A[0 : test split]
               bTrain = b[0 : test split]
               AValid = A[test split : test split + valid split]
               bValid = b[test split : test split + valid split]
               ATest = A[n - train split :]
               bTest = b[n - train split :]
               return ATrain, AValid, ATest, bTrain, bValid, bTest
           def toSparse(A, cols):
               row = []
               col = []
               data = []
               for i in range(len(A)):
                   for key, val in A[i].items():
                        row.append(i)
                        col.append(key - 1)
                        data.append(val)
               return coo matrix((data, (row, col)), shape=(len(A), cols)).tocsr
           ()
           def bConverter(b):
               for i in range(len(b)):
                   if (b[i] == 2):
                       b[i] = -1
               return b
           A news20 = toSparse(A news20, 1355191)
           b news20 = np.array(b news20)
           b \text{ news} 20 = b \text{ news} 20.\text{reshape}((b \text{ news} 20.\text{shape}[0], 1))
           print(A news20.shape)
           ATrain news20, AValid news20, ATest news20, bTrain news20, bValid new
           s20, bTest news20 = splitData(A news20, b news20)
           print('News data set loaded')
```

(19996, 1355191) News data set loaded

```
In [883]:
          initParams = {
               'lambda_': 0.001,
               'epsilon': 1e-2,
               'max iter': 2000,
               'alpha0': 1,
               't': 1,
               'beta0': np.random.uniform(-0.5, 0.5),
               'theta': 0.5,
               'x0 news20': np.random.uniform(-1, 1, size = (ATrain news20.shape
           [1], 1)),
               'sigma': 1e-12,
               'delta': 1e-12
           }
           proxGradData = []
In [884]:
           accelProxGradData = []
```

Download the dataset news20.binary dataset:

https://www.csie.ntu.edu.tw/~cjlin/libsvmtools/datasets/binary.html#news20.binary (https://www.csie.ntu.edu.tw/~cjlin/libsvmtools/datasets/binary.html#news20.binary) Check Assignment 4 for instructions about loading this dataset.

Question 1

Solve the l1-regularized logistic regression problem

$$ext{minimize}_{x \in \mathbb{R}^d, eta \in \mathbb{R}} \ \lambda \|x\|_1 + rac{1}{n} \sum_{i=1}^n \log(1 + \exp(-b_i(a_i^T x + eta)))$$

using

- 1) Proximal gradient descent
- 2) Accelerated proximal gradient descent
- 3) Proximal coordinate descent
- 4) Accelerated proximal coordinate descent

Tune the model parameter λ and any parameters that the algorithms have. You only have to tune λ once and use the same λ for all algorithms. Use similar techniques to Assignment 4 to tune λ . Plot the objective function (y-axis) vs running time in sec (x-axis) for all algorithms in the same plot. Make sure that the plots are clean and use appropriate legends. Report your generalization error. This should be measured in the same way as Assignment 4.

For training split the data into training (the first 15996 datapoints), validation (the next 2000 datapoints) and testing (the next 2000 datapoints). Do not use the testing data to influence training in any way. This means that you should compute the generalization error only once when you finish your work with this part of the assignment.

```
def lineSearch(beta, xCurrent, gradF, A, b):
In [885]:
               lambda = initParams['lambda ']
               alpha = initParams['alpha0']
               theta = initParams['theta']
               logisticRegression = LogisticRegression()
               proxGradDesc = ProxGradDesc()
               \#Compute\ F(x(alpha)):
               x alpha = proxGradDesc.proxAG(alpha, xCurrent, gradF)
               F x alpha = logisticRegression.objFunc(x alpha, A, b, beta)
               \#Compute\ F(x\ k):
               F x k = logisticRegression.objFunc(xCurrent, A, b, beta)
               #Compute theta * (l(x k) - l(x(alpha))):
               l_x_k = F_x_k
               l x alpha = logisticRegression.gx(x alpha) + logisticRegression.f
          x(xCurrent, A, b, beta) + np.dot(gradF.T, x alpha - xCurrent)
               while (F \times alpha > F \times k - theta * (l \times k - l \times alpha)):
                   alpha /= 2
                   \#Compute\ F(x(alpha)):
                   x alpha = proxGradDesc.proxAG(alpha, xCurrent, gradF)
                   F x alpha = logisticRegression.objFunc(x alpha, A, b, beta)
                   l x alpha = logisticRegression.gx(x alpha) + logisticRegressi
          on.fx(xCurrent, A, b, beta) + np.dot(gradF.T, x alpha - xCurrent)
               return alpha
          def validationError(bTrue, bPredict):
               t = len(bTrue)
               totalError = 0
               for i in range (t):
                   totalError += np.abs(bPredict[i] - bTrue[i])
               return totalError / (2 * t)
          def calcAccuracy(bTrue, bPredict):
               return 1 - .5 * validationError(bTrue, bPredict)
```

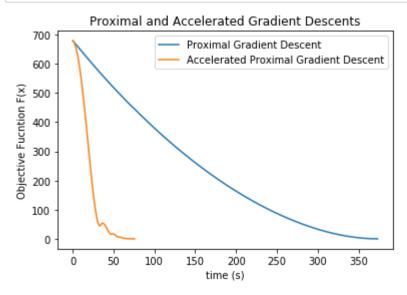
01/12/2019

In [886]: class LogisticRegression(object): def init (self): self.lambda = initParams["lambda "] def __str__(self): return 'L1-regularized Logistic Regression' **def** gx(self, x): lambda_ = self.lambda_ return lambda_ * np.linalg.norm(x, 1) def fx(self, x, A, b, beta): n = A.shape[0]aTx = A @ xsumLogistic = sum(np.log(1 + np.exp(-b * (aTx + beta))))return sumLogistic / n def objFunc(self, x, A, b, beta): g x = self.gx(x)f x = self.fx(x, A, b, beta) $return g_x + f_x$ def gradient(self, x, A, b, beta): n = A.shape[0]aTx = A @ xdenom = 1 / (1 + np.exp(b * (aTx + beta)))gradBeta = sum(-b * denom) / n gradFx = A.T @ (-b * denom)/nreturn gradFx, gradBeta def predict(self, x, A, beta): aTx = A @ xsigma = 1 + np.exp(-(aTx + beta))**return** 2 * (sigma < 2) - 1

```
In [887]: class ProxGradDesc():
               def __init__(self):
                   self.lambda = initParams['lambda ']
                   self.epsilon = initParams['epsilon']
                   self.beta = initParams['beta0']
                   self.logisticRegression = LogisticRegression()
               def str (self):
                   return 'Proximal Gradient Descent'
               def proxAG(self, alpha, x, gradF):
                   lambda_ = self.lambda_
                   xNext = np.zeros((x.shape[0], 1))
                   la = lambda_ * alpha
                   u = x - lambda * alpha * gradF
                   xNext[u >= la] = u[u >= la] - la
                   xNext[np.abs(u) \le la] = 0
                   xNext[u \leftarrow -la] = u[u \leftarrow -la] + la
                   return xNext
               def proxGrad(self, xCurrent, A, b):
                   beta = self.beta
                   max iters = initParams['max iter']
                   start = time.time()
                   for i in range(max iters):
                       gradF, gradB = self.logisticRegression.gradient(xCurrent,
          A, b, beta)
                       alpha = lineSearch(beta, xCurrent, gradF, A, b)
                       proxGradData.append([self.logisticRegression.objFunc(xCur
          rent, A, b, beta), time.time() - start])
                       if (i % 200 == 0):
                           print('Iteration:', i, 'out of', max_iters, 'iteratio
          ns')
                       xNext = self.proxAG(alpha, xCurrent, gradF)
                       Gx = (xCurrent - xNext) / alpha
                       if (np.linalg.norm(Gx) < self.epsilon):</pre>
                           return xNext, beta
                       else:
                           xCurrent = xNext
                           beta -= alpha * gradB
                   return xNext, beta
```

```
In [888]: | class AccelProxGradDesc():
              def __init__(self, lambda_ = initParams['lambda_']):
                   self.lambda = lambda
                   self.epsilon = initParams['epsilon']
                   self.beta = initParams['beta0']
                   self.t = initParams['t']
                   self.logisticRegression = LogisticRegression()
              def __str__(self):
                  return 'Accelerated Proximal Gradient Descent'
              def accelProxAG(self, alpha, y, gradF):
                   lambda = self.lambda
                   xNext = np.zeros((v.shape[0], 1))
                   la = lambda * alpha
                   u = y - lambda * alpha * gradF
                  xNext[u >= la] = u[u >= la] - la
                  xNext[np.abs(u) \le la] = 0
                  xNext[u \le -la] = u[u \le -la] + la
                   return xNext
              def accelProxGrad(self, xCurrent, A, b):
                   beta = self.beta
                  max iters = initParams['max iter']
                   tCurrent = self.t
                  yCurrent = xCurrent
                   start = time.time()
                   for i in range(max iters):
                       gradF, gradB = self.logisticRegression.gradient(xCurrent,
          A, b, beta)
                       alpha = lineSearch(beta, xCurrent, gradF, A, b)
                       accelProxGradData.append([self.logisticRegression.objFunc
          (xCurrent, A, b, beta), time.time() - start])
                       if (i \% 200 == 0):
                           print('Iteration:', i, 'out of', max iters, 'iteratio
          ns')
                       xNext = self.accelProxAG(alpha, yCurrent, gradF)
                       tNext = (1 + np.sqrt(1 + 4 * tCurrent ** 2)) / 2
                       vCurrent = xNext + ((tCurrent - 1)/tNext) * (xNext - xCur
          rent)
                       Gx = (xCurrent - xNext) / alpha
                       if (np.linalg.norm(Gx) < self.epsilon):</pre>
                           return xNext, beta
                       else:
                           xCurrent = xNext
                           tCurrent = tNext
                           beta -= alpha * gradB
                   return xNext, beta
```

```
In [889]:
          def calcError(x, beta):
              logisticRegression = LogisticRegression()
              bPredict = logisticRegression.predict(x, AValid news20, beta)
              error = validationError(bValid news20, bPredict)
              accuracy = calcAccuracy(bValid news20, bPredict)
              return error, accuracy
          x0 = initParams['x0 news20']
In [890]:
          print('Starting', ProxGradDesc(). str ())
          start = time.time()
          xProxGrad, betaProxGrad = ProxGradDesc().proxGrad(x0, ATrain news20,
          bTrain news20)
          end = time.time() - start
          print('Done', ProxGradDesc(). str (), 'in', end, 's')
          error, accuracy = calcError(xProxGrad, betaProxGrad)
          print('Error:', error, 'Accuracy:', accuracy)
          print()
          print('Starting', AccelProxGradDesc(). str ())
          start = time.time()
          xAccelProxGrad, betaAccelProxGrad = AccelProxGradDesc().accelProxGrad
          (x0, ATrain news20, bTrain news20)
          end = time.time() - start
          print('Done', AccelProxGradDesc(). str (), 'in', end, 's')
          error, accuracy = calcError(xAccelProxGrad, betaAccelProxGrad)
          print('Error:', error, 'Accuracy:', accuracy)
          Starting Proximal Gradient Descent
          Iteration: 0 out of 2000 iterations
          Iteration: 200 out of 2000 iterations
          Iteration: 400 out of 2000 iterations
          Iteration: 600 out of 2000 iterations
          Iteration: 800 out of 2000 iterations
          Iteration: 1000 out of 2000 iterations
          Done Proximal Gradient Descent in 373.2762362957001 s
          Error: [1.] Accuracy: [0.5]
          Starting Accelerated Proximal Gradient Descent
          Iteration: 0 out of 2000 iterations
          Done Accelerated Proximal Gradient Descent in 75.85138416290283 s
          Error: [1.] Accuracy: [0.5]
```



Did not have time to do Coordinate Descent

Question 2

Pick one algorithm that you want and plot the number of nonzeros in the output of the algorithm against λ . Explain in a couple of sentences (or less) what do you observe.

```
In [842]: | class NumNonZero():
              def __init__(self, lambda_):
                   self.lambda = lambda
                   self.epsilon = initParams['epsilon']
                   self.beta = initParams['beta0']
                   self.t = initParams['t']
                   self.logisticRegression = LogisticRegression()
              def __str__(self):
                  return 'Accelerated Proximal Gradient Descent'
              def accelProxAG(self, alpha, y, gradF):
                  lambda_ = self.lambda_
                  xNext = np.zeros((y.shape[0], 1))
                   la = lambda * alpha
                   u = y - lambda * alpha * gradF
                  xNext[u >= la] = u[u >= la] - la
                  xNext[np.abs(u) \le la] = 0
                  xNext[u \le -la] = u[u \le -la] + la
                   return xNext
              def accelProxGrad(self, xCurrent, A, b):
                   beta = self.beta
                  \max iters = 200
                   tCurrent = self.t
                  yCurrent = xCurrent
                   start = time.time()
                   for i in range(max iters):
                       gradF, gradB = self.logisticRegression.gradient(xCurrent,
          A, b, beta)
                       alpha = lineSearch(beta, xCurrent, gradF, A, b)
                       xNext = self.accelProxAG(alpha, yCurrent, gradF)
                       tNext = (1 + np.sqrt(1 + 4 * tCurrent ** 2)) / 2
                       vCurrent = xNext + ((tCurrent - 1)/tNext) * (xNext - xCur
          rent)
                       Gx = (xCurrent - xNext) / alpha
                       if (np.linalg.norm(Gx) < self.epsilon):</pre>
                           return xNext, beta
                       else:
                           xCurrent = xNext
                           tCurrent = tNext
                           beta -= alpha * gradB
                   return xNext, beta
```

```
In [859]: lambda list = np.power(10, np.linspace(-6, 0, 30))
          numNon0X = np.zeros like(lambda list)
          for i in range(len(lambda list)):
              numNonZero = NumNonZero(lambda list[i])
              print('Lambda =', lambda list[i])
              print('Starting', numNonZero.__str__(), 'number', i, 'of', len(la
          mbda_list))
              start = time.time()
              xAccelProxGrad, betaAccelProxGrad = numNonZero.accelProxGrad(x0,
          ATrain news20, bTrain news20)
              end = time.time() - start
              numNon0X[i] = np.count nonzero(xAccelProxGrad)
              print('Done', numNonZero.__str__(), 'number', i, 'of', len(lambda
          list), 'in', end, 's')
              print('Number of non zeros:', np.count nonzero(xAccelProxGrad))
              error, accuracy = calcError(xAccelProxGrad, betaAccelProxGrad)
              print('Error:', error, 'Accuracy:', accuracy)
              print()
```

Lambda = 1e-06

Starting Accelerated Proximal Gradient Descent number 0 of 30 Done Accelerated Proximal Gradient Descent number 0 of 30 in 0.384655 237197876 s

Number of non zeros: 1355187 Error: [0.528] Accuracy: [0.736]

Lambda = 1.6102620275609392e-06

Starting Accelerated Proximal Gradient Descent number 1 of 30 Done Accelerated Proximal Gradient Descent number 1 of 30 in 0.354690 0749206543 s

Number of non zeros: 1355187 Error: [0.528] Accuracy: [0.736]

Lambda = 2.592943797404667e-06

Starting Accelerated Proximal Gradient Descent number 2 of 30 Done Accelerated Proximal Gradient Descent number 2 of 30 in 0.356843 9483642578 s

Number of non zeros: 1355184 Error: [0.528] Accuracy: [0.736]

Lambda = 4.1753189365604e-06

Starting Accelerated Proximal Gradient Descent number 3 of 30 Done Accelerated Proximal Gradient Descent number 3 of 30 in 0.367183 4468841553 s

Number of non zeros: 1355180 Error: [0.528] Accuracy: [0.736]

Lambda = 6.723357536499335e-06

Starting Accelerated Proximal Gradient Descent number 4 of 30 Done Accelerated Proximal Gradient Descent number 4 of 30 in 0.354260 4446411133 s

Number of non zeros: 1355177 Error: [0.528] Accuracy: [0.736]

Lambda = 1.082636733874054e-05

Starting Accelerated Proximal Gradient Descent number 5 of 30 Done Accelerated Proximal Gradient Descent number 5 of 30 in 70.87307 31010437 s

Number of non zeros: 1338571 Error: [0.835] Accuracy: [0.5825]

Lambda = 1.7433288221999873e-05

Starting Accelerated Proximal Gradient Descent number 6 of 30 Done Accelerated Proximal Gradient Descent number 6 of 30 in 71.22521 686553955 s

Number of non zeros: 1328280

Error: [0.8445] Accuracy: [0.57775]

Lambda = 2.8072162039411757e-05

Starting Accelerated Proximal Gradient Descent number 7 of 30 Done Accelerated Proximal Gradient Descent number 7 of 30 in 73.61372 73311615 s

Number of non zeros: 1311522

Error: [0.8725] Accuracy: [0.56375]

Lambda = 4.520353656360241e-05

Starting Accelerated Proximal Gradient Descent number 8 of 30 Done Accelerated Proximal Gradient Descent number 8 of 30 in 74.36868 786811829 s

Number of non zeros: 1284874

Error: [0.9135] Accuracy: [0.54325]

Lambda = 7.278953843983146e-05

Starting Accelerated Proximal Gradient Descent number 9 of 30 Done Accelerated Proximal Gradient Descent number 9 of 30 in 73.06985 092163086 s

Number of non zeros: 1242130

Error: [0.9655] Accuracy: [0.51725]

Lambda = 0.00011721022975334806

Starting Accelerated Proximal Gradient Descent number 10 of 30 Done Accelerated Proximal Gradient Descent number 10 of 30 in 70.7144 0815925598 s

Number of non zeros: 1173004 Error: [1.] Accuracy: [0.5]

Lambda = 0.00018873918221350977

Starting Accelerated Proximal Gradient Descent number 11 of 30 Done Accelerated Proximal Gradient Descent number 11 of 30 in 70.7552 11353302 s

Number of non zeros: 1062055 Error: [1.] Accuracy: [0.5]

Lambda = 0.0003039195382313198

Starting Accelerated Proximal Gradient Descent number 12 of 30 Done Accelerated Proximal Gradient Descent number 12 of 30 in 70.5761 3897323608 s

Number of non zeros: 883938 Error: [1.] Accuracy: [0.5]

Lambda = 0.0004893900918477494

Starting Accelerated Proximal Gradient Descent number 13 of 30 Done Accelerated Proximal Gradient Descent number 13 of 30 in 70.4008 5554122925 s

Number of non zeros: 597989 Error: [1.] Accuracy: [0.5]

Lambda = 0.0007880462815669912

Starting Accelerated Proximal Gradient Descent number 14 of 30 Done Accelerated Proximal Gradient Descent number 14 of 30 in 68.6278 0714035034 s

Number of non zeros: 139617 Error: [1.] Accuracy: [0.5]

Lambda = 0.0012689610031679222

Starting Accelerated Proximal Gradient Descent number 15 of 30 Done Accelerated Proximal Gradient Descent number 15 of 30 in 57.0101 9215583801 s

Number of non zeros: 263 Error: [1.] Accuracy: [0.5]

Lambda = 0.0020433597178569417

Starting Accelerated Proximal Gradient Descent number 16 of 30

Done Accelerated Proximal Gradient Descent number 16 of 30 in 43.5631

00814819336 s

Number of non zeros: 12 Error: [1.] Accuracy: [0.5]

Lambda = 0.0032903445623126675

Starting Accelerated Proximal Gradient Descent number 17 of 30

Done Accelerated Proximal Gradient Descent number 17 of 30 in 30.7975

84533691406 s

Number of non zeros: 0 Error: [1.] Accuracy: [0.5]

Lambda = 0.005298316906283708

Starting Accelerated Proximal Gradient Descent number 18 of 30

Done Accelerated Proximal Gradient Descent number 18 of 30 in 22.7023

16761016846 s

Number of non zeros: 0

Error: [1.] Accuracy: [0.5]

Lambda = 0.008531678524172805

Starting Accelerated Proximal Gradient Descent number 19 of 30

Done Accelerated Proximal Gradient Descent number 19 of 30 in 15.9521 72756195068 s

Number of non zeros: 0

Error: [1.] Accuracy: [0.5]

Lambda = 0.013738237958832637

Starting Accelerated Proximal Gradient Descent number 20 of 30 Done Accelerated Proximal Gradient Descent number 20 of 30 in 12.1214

38980102539 s

Number of non zeros: 0

Error: [1.] Accuracy: [0.5]

Lambda = 0.022122162910704502

Starting Accelerated Proximal Gradient Descent number 21 of 30

Done Accelerated Proximal Gradient Descent number 21 of 30 in 8.55515 9091949463 s

Number of non zeros: 0

Error: [1.] Accuracy: [0.5]

Lambda = 0.03562247890262444

Starting Accelerated Proximal Gradient Descent number 22 of 30

Done Accelerated Proximal Gradient Descent number 22 of 30 in 6.21954

607963562 s

Number of non zeros: 0

Error: [1.] Accuracy: [0.5]

Lambda = 0.05736152510448681

Starting Accelerated Proximal Gradient Descent number 23 of 30

Done Accelerated Proximal Gradient Descent number 23 of 30 in 4.27811

9087219238 s

Number of non zeros: 0

Error: [1.] Accuracy: [0.5]

Lambda = 0.09236708571873865

Starting Accelerated Proximal Gradient Descent number 24 of 30

Done Accelerated Proximal Gradient Descent number 24 of 30 in 3.21043

4675216675 s

Number of non zeros: 0 Error: [1.] Accuracy: [0.5]

Lambda = 0.14873521072935117

Starting Accelerated Proximal Gradient Descent number 25 of 30 Done Accelerated Proximal Gradient Descent number 25 of 30 in 2.51329

56504821777 s

Number of non zeros: 0 Error: [1.] Accuracy: [0.5]

Lambda = 0.2395026619987486

Starting Accelerated Proximal Gradient Descent number 26 of 30 Done Accelerated Proximal Gradient Descent number 26 of 30 in 1.59848 0224609375 s

Number of non zeros: 0 Error: [1.] Accuracy: [0.5]

Lambda = 0.38566204211634725

Starting Accelerated Proximal Gradient Descent number 27 of 30 Done Accelerated Proximal Gradient Descent number 27 of 30 in 1.25519 82402801514 s

Number of non zeros: 0 Error: [1.] Accuracy: [0.5]

Lambda = 0.6210169418915616

Starting Accelerated Proximal Gradient Descent number 28 of 30 Done Accelerated Proximal Gradient Descent number 28 of 30 in 0.92773 4375 s

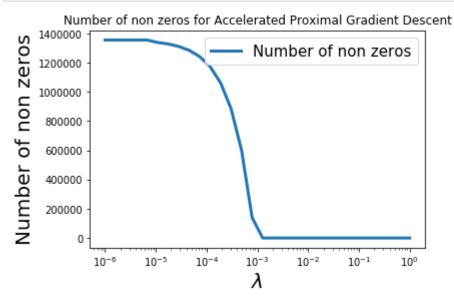
Number of non zeros: 0 Error: [1.] Accuracy: [0.5]

Lambda = 1.0

Starting Accelerated Proximal Gradient Descent number 29 of 30 Done Accelerated Proximal Gradient Descent number 29 of 30 in 0.59047 38903045654 s

Number of non zeros: 0 Error: [1.] Accuracy: [0.5]

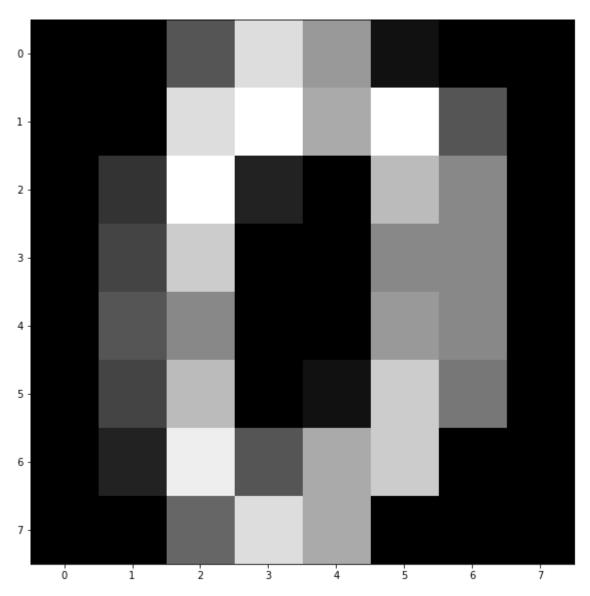
```
In [860]: plt.plot(lambda_list, numNon0X, label = "Number of non zeros", linewi
    dth = 3)
    plt.legend(fontsize = 15)
    plt.xlabel("$\lambda$", fontsize = 20)
    plt.ylabel("Number of non zeros", fontsize = 20)
    plt.xscale('log')
    plt.title('Number of non zeros for Accelerated Proximal Gradient Desc
    ent')
    plt.show()
```



From the graph and the printed data, we can see that the number of non zero outputs goes to zero as lambda goes to increases. This is expected since lambda is the weight for the regulization term. As we are minimizing the objective function, the regulization term will be at its lowest when x is all zeros. So if we increase lambda, the minimum of the function will most likely be found when x is all zeros.

Label Propagation

Load datasets from sklearn In [892]: from sklearn import datasets # Load the digits dataset. # See here for a distription of this dataset: # https://scikit-learn.org/stable/modules/generated/sklearn.datasets. load digits.html#sklearn.datasets.load digits digits = datasets.load digits() # Briefly, this dataset has 1797 images of digits. The digits are fro m 0 to 9. # There are about 180 images per digit. # digits.data includes the vectorized images of the digits # digits.target includes the labels of the images # digits.images includes the images in non-vectorized format # Example: number 0 import matplotlib.pyplot as plt plt.figure(1, figsize=(10, 10)) plt.imshow(digits.images[0], cmap='gray') plt.show()



Let x_i and xj be two vectorized images. Compute their pairwise similarity by \$\$ w{ij} = \exp(-\gamma |x_i-x_j|_2^2), \$\$ where γ is parameter tha controls the importance of the distance $|x_i-x_j|_2^2$. If $\gamma=1/(2\sigma^2)$ then this is called the Gaussian kernel and σ^2 represents the variance. You can compute the similarities by using: $\frac{\text{https://scikit-learn.org/stable/modules/generated/sklearn.metrics.pairwise.rbf_kernel.html (https://scikit-learn.org/stable/modules/generated/sklearn.metrics.pairwise.rbf_kernel.html)}.$ Alterantively, you can compute the pairwise similarities manually, by using the above formula.

Sparsify the graph by setting small values to zero. The thresholding tolerance is a parameter that you will have to tune. Generally, there is a consensus that sparse similarity matrices result in better performance. So, you should expect that, but have in mind that this is only an empirical observation.

```
In [893]: # Example
# Pairwise weights matrix
W = sklearn.metrics.pairwise.rbf_kernel(digits.data)

# Remove self-similarity
W = W - np.eye(len(W))

# Sparsify (tune the tolerance parameter for your problem)
tol = le-9
W[abs(W) < tol] = 0.0</pre>
```

This dataset provides labels for every image of a digit. However, in this assignment we want to check how the label propagation model works. Use all images, i.e., 1797 images of digits, but use only 10 labels per type of digit, e.g., use the first 10 labels for digit 0, the first 10 labels of digit 1 etc. This means that you should use 100 labels in total out of 1797. Do not use the remaining labels for training the model. The remaining labels will be used only to measure accuracy of the model.

```
In [894]:
          # digits.data includes the vectorized images of the digits
          # digits.target includes the labels of the images
          # digits.images includes the images in non-vectorized format
          def createDataSet(digits, size, numbers = 10):
              knownSetLabel = np.array([])
              knownSetInd = np.array([])
              unknownSetInd = np.array([])
              for i in range(numbers):
                  knownSetInd = np.append(knownSetInd, np.where(digits.target =
          = i)[0][: size])
                  unknownSetInd = np.append(unknownSetInd, np.where(digits.targ
          et == i)[0][size :])
              knownSetLabel = np.append(knownSetLabel, digits.target[knownSetIn
          d.astype(int)])
              return knownSetInd, knownSetLabel, unknownSetInd
```

Question 3

Train the label propagation model

$$ext{minimize}_{b \in \mathbb{R}^n} \; rac{1}{2} \sum_{i=1}^n \sum_{j=1}^n w_{ij} (b_i - b_j)^2 + \sum_{i=1}^{m-n} \sum_{j=1}^n ar{w}_{ij} (ar{b}_i - b_j)^2$$

using the coordinate method that was discussed in Lecture 14 for the label propagation problem. Make sure that you understand the derivation of the algorithm before you apply it to train the label propagation model.

The output of the algorithm should be the soft labels b for each image. However, these soft labels are not going to be integers (actual labels). Therefore, we have to round the output labels b to get integer labels for each image. One possible way to round the output is to simply find the integer label (possible options 0 to 9) that a soft label is closest to. Then pick this label as the label of the image.

After you label all images, use the function in https://scikit-

<u>learn.org/stable/modules/generated/sklearn.metrics.precision_recall_fscore_support.html</u> (https://scikit-learn.org/stable/modules/generated/sklearn.metrics.precision_recall_fscore_support.html) to measure precision, recall and f1_score. You only have to provide the true and the predicted labels when using this function. Do not change the default parameters.

If you want to learn more about precision, recall and f1_score check Wikipedia: precision and recall: https://en.wikipedia.org/wiki/Precision_and_recall (https://en.wikipedia.org/wiki/Precision_and_recall) f1_score: https://en.wikipedia.org/wiki/F1_score (https://en.wikipedia.org/wiki/F1_score).

Report precision, recall and f1score of your method.

```
In [896]: import pprint
    bBarI, bBar, bI = createDataSet(digits, 10)
    b = labelProp(bBar, bBarI, bI, W)

precision, recall, f1_score, support = sklearn.metrics.precision_recall_fscore_support(digits.target, b)
    pd.DataFrame(data = {'Precision': precision, 'Recall': recall, 'f1 Score': f1_score, 'Num of Labels': support})
```

Out[896]:

	Precision	Recall	f1 Score	Num of Labels
0	1.000000	0.994382	0.997183	178
1	0.960000	0.263736	0.413793	182
2	0.593750	0.107345	0.181818	177
3	0.430341	0.759563	0.549407	183
4	0.729167	0.966851	0.831354	181
5	0.541254	0.901099	0.676289	182
6	0.648746	1.000000	0.786957	181
7	0.491124	0.927374	0.642166	179
8	0.466667	0.120690	0.191781	174
9	1.000000	0.055556	0.105263	180

Question 4

Increase the number of labels that are used in the model from 10 per digit to 20 per digit. Re-train and report precision, recall and f1score of your method. Explain in a couple of sentences what do you observe.

```
In [897]: bBarI, bBar, bI = createDataSet(digits, 20)
b = labelProp(bBar, bBarI, bI, W)

precision, recall, f1_score, support = sklearn.metrics.precision_recall_fscore_support(digits.target, b)
pd.DataFrame(data = {'Precision': precision, 'Recall': recall, 'f1 Score': f1_score, 'Num of Labels': support})
```

Out[897]:

		Precision	Recall	f1 Score	Num of Labels
_	0	1.000000	0.994382	0.997183	178
	1	0.953125	0.335165	0.495935	182
	2	0.745614	0.960452	0.839506	177
	3	0.659898	0.710383	0.684211	183
	4	0.660377	0.966851	0.784753	181
	5	0.826389	0.653846	0.730061	182
	6	0.628472	1.000000	0.771855	181
	7	0.666667	0.927374	0.775701	179
	8	0.328000	0.235632	0.274247	174
	9	0.950000	0.316667	0.475000	180

Local Graph Clustering

```
In [898]: import numpy as np
          # Import networkx to load the data
          import networkx as nx
          # Provide the path of the dataset JohnsHopkins55
          path = 'JohnsHopkins55.graphml'
          # This dataset is a Facebook anonymized dataset on a particular day i
          n September 2005
          # for a student social network at John Hopkins university. The graph
           is unweighted
          # and it represents ``friendship'' ties. This graph has 5157 nodes an
          d 186572 edges.
          # This dataset comes along with 6 features, i.e., second major, high
           school, gender, dorm, major index and year.
          # We construct ``ground truth'' clusters by using the features for ea
          ch node.
          # In particular, we consider nodes with the same value of a feature t
          o be a cluster, e.g., students of year 2009.
          # For this assignment we will attempt to find the students of year 20
          09 using local graph clustering.
          # Load the data as a graph
          G = nx.read graphml(path).to undirected()
          # Construct the adjacency matrix
          A = nx.adjacency matrix(G).astype(np.float64)
          ALocalGraphCluster = A
          # Indices of nodes (students) of class year 2009
          # There should be 910 students of class year 2009.
          # We will use these students to measure performance of the model
          students of year 2009 = []
          ct = 0
          for node in G.nodes(data=True):
              if node[1]['year'] == 2009:
                  students_of_year_2009.append(ct)
              ct += 1
          print(len(students of year 2009))
```

910

The target cluster is the students of year 2009 in the list: students_of_year_2009.

For this problem you will have to use the coordinate algorithm that was introduced in Lecture 14.

The coordinate algorithm returns a vector p. To get an actual cluster from vector p you will have to round the output of the coordinate algorithm using the rounding technique that was discussed in Lecture 14.

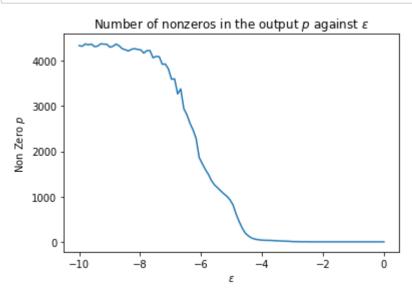
To measure accuracy of the rounded output you can use number of true positives and the number of false positives of the predicted cluster. True positives are the predicted nodes that are also part of class year 2009. False positive are the number of predicted nodes that are not part of the class year 2009.

```
In [899]:
          #A: Adjacency Martix
          #degree: array of degrees of each nodes
          def localGraphCluster(A, degree, seed, alpha, e, max iters = 10000):
              n = A.shape[0]
              p = np.zeros(n)
               r = np.zeros(n)
              nodes = {seed}
               r[seed] = 1
              i = 0
              while i < max iters and len(nodes) > 0:
                   u = random.sample(nodes, 1)[0]
                   p[u] = p[u] + alpha * r[u]
                   v = A[u].indices
                   r[v] += (0.5 * (1 - alpha) * r[u]) / degree[u]
                   r[u] *= 0.5 * (1 - alpha)
                  if r[u] < e * degree[u]:
                       nodes.remove(u)
                   nodes.update({i for i in v if r[i] >= e * degree[i]})
                   i += 1
               return p
          def getCluster(p, A, degree):
              non0Nodes = np.argwhere(p > 0).flatten()
              non0Nodes = non0Nodes[np.argsort(-p[non0Nodes]/degree[non0Nodes
          ])]
              edges = np.zeros like(p)
              degreeSum = 0
              conductances = []
              for i, node in enumerate(non0Nodes):
                   neighbours = A[node].indices
                   degreeSum += len(neighbours)
                   edges[neighbours] = 1
                   edges[non0Nodes[: i + 1]] = 0
                   conductances.append(sum(edges) / degreeSum)
               cluster = non0Nodes[: np.argmin(conductances) + 1]
               return cluster
```

Question 5

1) Report the number of nonzero nodes in the output of the coordinate algorithm (vector p) as the tolerance parameter ϵ goes to zero. In particular, plot the number of nonzeros in the output p against ϵ . The number of nonzeros should increase as ϵ goes to zero.

```
In [900]:
          seed = 5000
          alpha = 0.1
          non0P = []
          A = ALocalGraphCluster
          #epsilon from 1e-10 to 1
          epsilonList = np.power(10, np.linspace(0, -10, 100))
          degree = np.array([A[i].sum() for i in range(A.shape[0])], dtype = in
          t)
          for epsilon in epsilonList:
              p = localGraphCluster(A, degree, seed, alpha, epsilon)
              non OP.append(p[p > 0].size)
          plt.figure()
          plt.plot(np.log10(epsilonList), non0P)
          plt.xlabel('$\epsilon$')
          plt.ylabel('Non Zero $p$')
          plt.title('Number of nonzeros in the output $p$ against $\epsilon$')
          plt.show()
```



Question 6

2) Run the coordinate algorithm using as seed node the node with index 5000. Round the output to get an actual cluster back. Measure true positives and false positives of the predicted cluster. Report these two numbers.

```
In [901]:
          epsilon = 1e-5
          seed5000 = 5000
          alpha = 0.1
          p5000 = localGraphCluster(A, degree, seed5000, alpha, epsilon)
          cluster = getCluster(p5000, A, degree)
          truePos = 0
          falsePos = 0
          for i in cluster:
              if i in students_of_year_2009:
                   truePos += 1
              else:
                   falsePos += 1
          print('True and False positives for seed:', seed5000)
          print('True Positive:', truePos)
          print('False Positive:', falsePos)
```

True and False positives for seed: 5000 True Positive: 744 False Positive: 71

Question 7

3) Run the coordinate algorithm using as seed node the node with index 146. Round the output to get an actual cluster back. Measure true positives and false positive of the predicted cluster. Report these two numbers. How do they compare to the results in 2)? Are they different? If yes, why do you think this might have happened? Explain intuitively in a couple of sentences why you think this might happen.

Marks: 5

True and False positives for seed: 146 True Positive: 572 False Positive: 111

Compare nodes 5000 and 146

Yes the results for Seed node 5000 and Seed node 146 are indeed different. The number of true positives decrease, and we get more False Positives when we start with node 146. This is because node 146 only has 3 neighbours, thus we miss out on more classes

Reccomender System

Download the data matrix A from the webiste https://grouplens.org/(https://grouplens.org/). Navigate to the dataset tab https://grouplens.org/datasets/movielens/) and download the dataset MovieLens 100K. Read the description of the dataset carefully. This dataset includes 5 data matrices A. We only need to use one of these in this assignment. In particular, use the dataset u1.base. Use the corresponding testing data u1.test to measure the accuracy of your model. For this question only it's ok if you use the testing data to tune parameters. Ignore the timestap feature in the dataset. We do not need this for this problem.

```
train = np.loadtxt('ml-100k/ul.base')[:, :3]
In [903]:
          test = np.loadtxt('ml-100k/u1.test')[:, :3]
          def getA(data, n, m):
              data[:, :2] = data[:, :2] - 1
              A = np.zeros((n, m))
              for i in range(len(data)):
                   user = int(data[i][0])
                  movie = int(data[i][1])
                   rating = data[i][2]
                   A[user][movie] = rating
               return A
          ATrainRS = getA(train, 943, 1682)
          ATestRS = getA(test, 943, 1682)
          x0 = np.random.randint(1, 6, ATrainRS.shape)
          knownValues = ATrainRS != 0
          knownTestValues = ATestRS != 0
```

Question 8

Solve the penalized problem

$$ext{minimize } \lambda \|X\|_* + \sum_{ij \, | \, A_{ij} \, is \, known} (X_{ij} - A_{ij})^2,$$

where $||X||_*$ is the nuclear norm, using proximal gradient descent. This algorithm requires thresholding the singular values of X at each iteration. You can use

https://docs.scipy.org/doc/scipy/reference/generated/scipy.linalg.svd.html

(https://docs.scipy.org/doc/scipy/reference/generated/scipy.linalg.svd.html) (sparse matrices) or

https://docs.scipy.org/doc/numpy/reference/generated/numpy.linalg.svd.html

(https://docs.scipy.org/doc/numpy/reference/generated/numpy.linalg.svd.html) (dense matrices).

To check your solution against u1.test, report your score as follows. Let X be the output of the proximal gradient method, let A' and Ω' be the data matrix and the non-zero entry list for the test data. Then the score is

$$rac{1}{|\Omega'|} \sum_{(i,j) \in \Omega'} |X_{ij} - A'_{ij}|.$$

In other words, the average absolute deviation of the computed ratings from the predicted ratings. Report your score.

```
In [904]: def prox(lambda , alpha, xCurrent, gradF):
               z = xCurrent - alpha * gradF
               u, s, vT = np.linalg.svd(z)
               s = np.maximum(s - alpha * lambda_, 0)
               sDense = np.zeros((u.shape[1], vT.shape[0]))
               sDense[:len(s), :len(s)] = np.diag(s)
               xNext = (u.dot(sDense)).dot(vT)
               return xNext, s
           def lineSearchRS(xCurrent, A, gradF):
               alpha = initParams['alpha0']
               theta = initParams['theta']
               lambda = 50
               rs = ReccomenderSystem()
               \#F(x | k)
               s = np.linalg.svd(xCurrent, compute uv = False)
               F x k = rs.objFunc(xCurrent, A, s)
               \#l(x | k)
               l_x_k = F_x_k
               #x(alpha)
               x_alpha, s = prox(lambda_, alpha, xCurrent, gradF)
               #F(x(alpha))
               F_x_alpha = rs.objFunc(x_alpha, A, s)
               \#l(x(alpha))
               l \times alpha = rs.gx(x alpha, s) + rs.fx(xCurrent, A) + (gradF * (x))
           alpha - xCurrent)).sum()
               while (F \times alpha > F_x_k - theta * (l_x_k - l_x_alpha)):
                   alpha /= 2
                   #x(alpha)
                   x_alpha, s = prox(lambda_, alpha, xCurrent, gradF)
                   \#F(x(alpha))
                   F \times alpha = rs.objFunc(x alpha, A, s)
                   \#l(x(alpha))
                   l_x_alpha = rs.gx(x_alpha, s) + rs.fx(xCurrent, A) + (gradF *
           (x alpha - xCurrent)).sum()
               return alpha
           def error(x, ATest):
               x = np.round(x)
               error = np.linalg.norm(x[knownTestValues] - ATest[knownTestValues
           1, 1)/knownTestValues.sum()
               return error
```

```
In [905]: class ReccomenderSystem():
               def init (self):
                   self.lambda = 50 #initParams["lambda "]
               def __str__(self):
                   return 'Reccomender System'
               def gx(self, X, s):
                   lambda = self.lambda
                   return lambda_ * np.sum(s)
               def fx(self, X, A):
                   return np.sum((X[knownValues] - A[knownValues]) ** 2)
               def objFunc(self, X, A, s):
                   g_x = self.gx(A, s)
                   f x = self.fx(X, A)
                   return g \times + f \times
               def gradient(self, X, A):
                   gradF = np.zeros like(A)
                   gradF[knownValues] = 2 * (X[knownValues] - A[knownValues])
                   return gradF
```

```
In [906]:
          def penalizer(xCurrent, A):
               rs = ReccomenderSystem()
               \max iters = 200
               lambda = 10
               for i in range(max_iters):
                   gradF = rs.gradient(xCurrent, A)
                   alpha = lineSearchRS(xCurrent, A, gradF)
                   xNext, s = prox(lambda , alpha, xCurrent, gradF)
                   if (i % 20 == 0):
                       print('Iteration:', i, 'out of', max_iters, 'iterations')
                       print('Objective Function:', rs.objFunc(xNext, A, s))
                  Gx = (xCurrent - xNext) / alpha
                   if (np.linalg.norm(Gx) < 1e-2):</pre>
                       return xNext
                   else:
                       xCurrent = xNext
               return xNext
```

```
rs = ReccomenderSystem()
In [667]:
          x0[knownValues] = ATrainRS[knownValues]
          x = penalizer(x0, ATrainRS)
          score = error(x, ATest)
          print('Score:', score)
```

Iteration: 0 out of 200 iterations Objective Function: 2224034.5471557504 Iteration: 20 out of 200 iterations Objective Function: 342708.27298988577 Iteration: 40 out of 200 iterations Objective Function: 332658.91948083654 Iteration: 60 out of 200 iterations Objective Function: 327585.0407599081 Iteration: 80 out of 200 iterations Objective Function: 324214.87051402545 Iteration: 100 out of 200 iterations Objective Function: 321716.4282558738 Iteration: 120 out of 200 iterations Objective Function: 319755.3231732809 Iteration: 140 out of 200 iterations Objective Function: 318164.5651491285 Iteration: 160 out of 200 iterations Objective Function: 316847.3353249789 Iteration: 180 out of 200 iterations Objective Function: 315742.0705507721

Score: 0.7265

Question 9

Now consider the factorized problem

$$ext{minimize}_{L,R} \ rac{\lambda}{2} ig(\|L\|_F^2 + \|R\|_F^2 ig) + \|W \circ (LR^T - A)\|_F^2$$

See Lecture 15 about details for this problem. Solve this problem and obtain the matrix $X=LR^T$. Use X as the predicted rating and report your score. For calculating derivatives of matrix functions you can have a look here: https://www.math.uwaterloo.ca/~hwolkowi/matrixcookbook.pdf (https://www.math.uwaterloo.ca/~hwolkowi/matrixcookbook.pdf).

```
In [907]: A = ATrainRS
          m, n = A.shape
          W = np.zeros like(A)
          W[knownValues] = 1
          k = 100
          L0 = np.random.rand(m, k)
          R0 = np.random.rand(n, k)
          class FactorizedReccomenderSystem():
              def init (self, lambda ):
                   self.lambda_ = lambda_
              def __str__(self):
                   return 'Factorized Reccomender System'
              def objFunc(self, L, R, W, A):
                   lambda = self.lambda
                   Lnorm = np.linalg.norm(L, ord = 'fro') ** 2
                   Rnorm = np.linalg.norm(R, ord = 'fro') ** 2
                   right = np.linalg.norm(W * (L @ R.T - A), ord = 'fro') ** 2
                   return (lambda_ / 2) * (Lnorm + Rnorm) + right
              def gradL(self, L, R, W, A):
                   lambda = self.lambda
                   return lambda * L + \frac{1}{2} * (W * (L @ R.T - A)) @ R
              def gradR(self, L, R, W, A):
                   lambda = self.lambda
                   return lambda_ * R + 2 * ((W * (L @ R.T - A)).T @ L)
              def LLineSearch(self, L, R, W, A):
                   lambda_ = self.lambda_
                   alpha = 1
                   f \times k = objFunc(lambda , L, R, W, A)
                   gL = gradL(lambda_, L, R, W, A)
                   LNext = L - alpha * gL
                   while(objFunc(lambda_, LNext, R, W, A) > f_x_k):
                       alpha /= 2
                       LNext = L - alpha * gL
                   return alpha
              def RLineSearch(self, L, R, W, A):
                   lambda = self.lambda_
                   alpha = 1
                   f \times k = objFunc(lambda , L, R, W, A)
                   gR = gradR(lambda, L, R, W, A)
                   RNext = R - alpha * gR
                   while(objFunc(lambda_, L, RNext, W, A) > f_x_k):
                       alpha /= 2
                       RNext = R - alpha * gR
                   return alpha
              def factorizedPenalizer(self, L, R, W, A):
                   lambda_ = self.lambda_
                   max iters = initParams['max iter']
                   for i in range(max iters):
                       #Fix L and solve wrt R
```

```
alphaR = RLineSearch(lambda_, L, R, W, A)
gR = gradR(lambda_, L, R, W, A)

R = R - alphaR * gR

#Fix R and solve wrt L
alphaL = LLineSearch(lambda_, L, R, W, A)
gL = gradL(lambda_, L, R, W, A)
L = L - alphaL * gL

if (i % 200 == 0):
    print('Iteration:', i, 'out of', max_iters, 'iterations')

print('Objective Function:', objFunc(lambda_, L, R, W, A))

return L, R
```

```
In [908]: lambda_ = 10
    factorizedReccomenderSystem = FactorizedReccomenderSystem(lambda_)
    L, R = factorizedReccomenderSystem.factorizedPenalizer(L0, R0, W, A)
    x = L @ R.T
    score = error(x, ATest)
    print('Score:', score)
```

Iteration: 0 out of 2000 iterations Objective Function: 10669077.629436143 Iteration: 200 out of 2000 iterations Objective Function: 81709.61436792536 Iteration: 400 out of 2000 iterations Objective Function: 72281.12131890278 Iteration: 600 out of 2000 iterations Objective Function: 70697.90595178634 Iteration: 800 out of 2000 iterations Objective Function: 70166.04706762842 Iteration: 1000 out of 2000 iterations Objective Function: 69894.50762842332 Iteration: 1200 out of 2000 iterations Objective Function: 69730.00617448597 Iteration: 1400 out of 2000 iterations Objective Function: 69620.10686353974 Iteration: 1600 out of 2000 iterations Objective Function: 69541.97711507631 Iteration: 1800 out of 2000 iterations Objective Function: 69484.5006241859 Score: 0.75265

Nonnegative Matrix Factorization: facial feature extraction

```
In [909]:
          # This piece of code is for loading data and visualizing
          # the first 6 images in the dataset.
          # Useful packages for loading the data and plotting
          from numpy.random import RandomState
          import matplotlib.pyplot as plt
          from sklearn.datasets import fetch olivetti faces
          n row, n col = 2, 3
          image shape = (64, 64)
          rng = RandomState(0)
          # Useful function for plotting
          def plot gallery(title, images, n col=n col, n row=n row, cmap=plt.cm
          .gray):
              plt.figure(figsize=(2. * n_col, 2.26 * n_row))
              plt.suptitle(title, size=16)
              for i, comp in enumerate(images):
                   plt.subplot(n_row, n_col, i + 1)
                   vmax = max(comp.max(), -comp.min())
                   plt.imshow(comp.reshape(image shape), cmap=cmap,
                              interpolation='nearest',
                              vmin=-vmax, vmax=vmax)
                   plt.xticks(())
                  plt.yticks(())
              plt.subplots adjust(0.12, 0.05, 0.99, 0.75, 0.04, 0.)
          # Load faces data
          dataset = fetch olivetti faces(shuffle=True, random state=rng)
          \# Store the vectorized images. Each image has dimensions 64 \times 64.
          faces = dataset.data
          print("Dataset consists of %d faces" % len(faces))
          plot gallery("First 6 faces from the dataset", faces[:6], 6, 1)
```

Dataset consists of 400 faces

First 6 faces from the dataset













```
In [910]: | class NonNegMatFac():
              def __init (self, data):
                   self.sigma = initParams['sigma']
                   self.delta = initParams['delta']
                   self.X = data.T
              def __str__(self):
                   return 'Non-negative Matrix Factorization'
              def objFunc(self, W, H):
                  X = self.X
                   return 0.5 * np.linalg.norm(W @ H - X, ord='fro') ** 2
              def gradH(self, W, H):
                  X = self.X
                   return W.T @ (W @ H - X)
              def gradW(self, W, H):
                   X = self.X
                   return (W @ H - X) @ H.T
              def HBar(self, W, H):
                   gradH = self.gradH(W, H)
                   HBar = np.where(gradH >= 0, H, np.maximum(H, self.sigma * np.
          ones((H.shape))))
                   return HBar
              def WBar(self, W, H):
                   gradW = self.gradW(W, H)
                   WBar = np.where(gradW >= 0, W, np.maximum(W, self.sigma * np.
          ones((W.shape))))
                   return WBar
              def bars(self, W, H):
                   HBar = self.HBar(W, H)
                   WBar = self.WBar(W, H)
                   return HBar, WBar
```

Question 10

Use the data in "faces" from the above code for the nonnegative matrix factorization problem

$$egin{aligned} ext{minimize}_{W,H} & \|WH - X\|_F^2 \ ext{subj. to } W_{ij} \geq 0 \ orall i,j \ H_{ij} \geq 0 \ orall i,j \end{aligned}$$

where $W \in \mathbb{R}^{m \times r}$ and $H \in \mathbb{R}^{r \times n}$. Set parameter r=6 in the nonnegative factorization problem. Extract the features matrix W by solving the non-negative matrix factorization problem. Report the 6 features of the faces dataset, i.e., the 6 columns of matrix W. You can report the features by visualizing them in a similar way to the above example.

```
In [911]: def nonNegMatFac(data, r = 6):
              m, n = data.T.shape
              max iters = initParams['max iter']
              nnmf = NonNegMatFac(data)
              W = np.random.uniform(0, 1, size = (m, r))
              H = np.random.uniform(0, 1, size = (r, n))
              delta = initParams['delta']
              for i in range(max iters):
                  HBar, WBar = nnmf.bars(W, H)
                  H = H - np.divide(HBar, ((W.T @ W @ HBar) + delta)) * nnmf.gr
          adH(W, H)
                  W = W - np.divide(WBar, ((WBar @ H @ HBar.T) + delta)) * nnmf
           .gradW(W, H)
                  #Normalize
                  S = np.eye(r) / np.sum(W, 0)
                  W = W \otimes S
                  H = np.linalg.inv(S) @ H
                  if (i \% 200 == 0):
                       print('Iteration:', i, 'out of', max_iters, 'iterations')
                      print('Objective Function:', nnmf.objFunc(W, H))
              return W, H
```

```
In [912]: W, H = nonNegMatFac(faces)
plot_gallery("Features", W.T[:6], 6, 1)
```

Iteration: 0 out of 2000 iterations Objective Function: 25793.681538200235 Iteration: 200 out of 2000 iterations Objective Function: 7010.241055904761 Iteration: 400 out of 2000 iterations Objective Function: 6910.017278901699 Iteration: 600 out of 2000 iterations Objective Function: 6891.487934288256 Iteration: 800 out of 2000 iterations Objective Function: 6882.712619091637 Iteration: 1000 out of 2000 iterations Objective Function: 6877.521901832734 Iteration: 1200 out of 2000 iterations Objective Function: 6873.825724841319 Iteration: 1400 out of 2000 iterations Objective Function: 6871.02090566685 Iteration: 1600 out of 2000 iterations Objective Function: 6868.694681355642 Iteration: 1800 out of 2000 iterations Objective Function: 6866.78572314461

Features













Question 11

Visualize the approximation of the first 6 images using the feature matrix W and the importance matrix H. Note that each face can be approaximated by a linear combination of features (columns) from matrix W. The positive coefficients of this linear combination are given by matrix H.

```
In [913]: WH = W @ H
   plot_gallery('Approximations of First 6 Images', WH.T[:6], 6, 1)
```

Approximations of First 6 Images













Question 12

Use matrix H to find the 3 most important features of the first 6 faces. To find the 3 most important features for an image pick the columns of W with the largest coefficients in matrix H.

```
In [914]: first6 = faces[: 6]
    top3Features = []
    for i in range(6):
        top3Features.append(np.argsort(-H[:,i])[:3])
    top3Features = np.asarray(top3Features)
    for i in range(3):
        first6 = np.append(first6, W.T[top3Features.T[i]], 0)
    for i in range(4):
        if (i == 0):
            plot_gallery('Original', first6[0: (i + 1) * 6], 6, 1)
        else:
            plot_gallery('Feature ' + str(i), first6[i * 6: (i + 1) * 6],
        6, 1)
```

