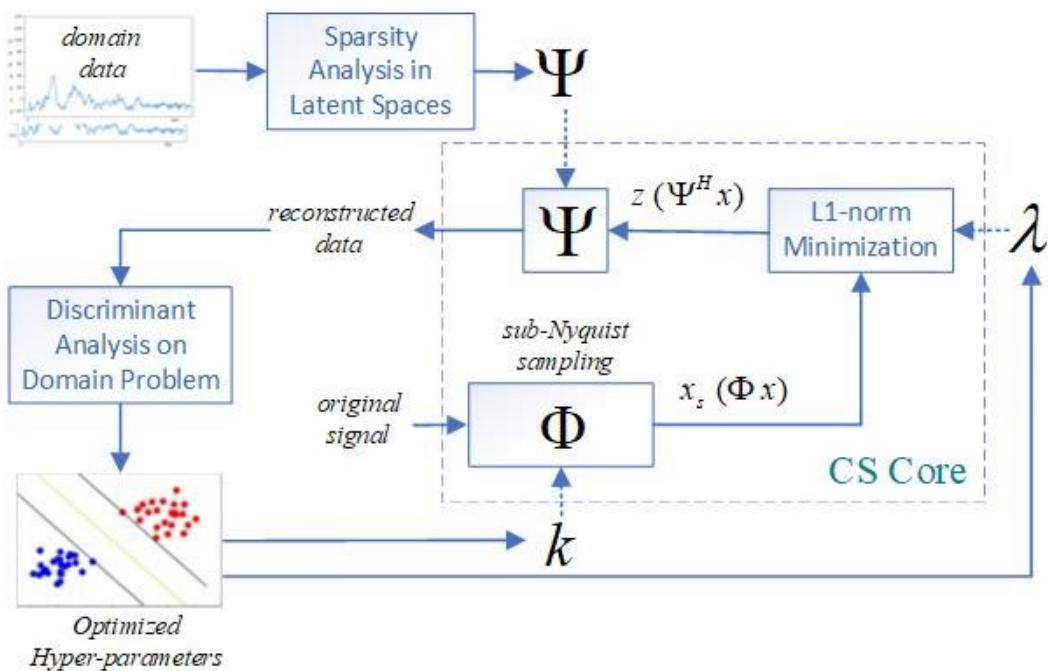


Adaptive Compressed Sensing of Raman Spectroscopic Profiling Data for Discriminative Tasks



Core Compressed Sensing (CS) Process

Sampling

$$x_s = \Phi x$$

Sparsity hypothesis

$$x = \Psi z$$

Compressed Sensing / sub-Nyquist Sampling

$$x_s = \Phi \Psi z$$

Reconstruction of latent z

$$\text{minimize } \|z\|_1 \quad s.t. \quad x_s = \Phi \Psi z$$

Inverse-transform to orginal space

$$x = \Psi z$$

Load Data

```
In [5]: import os
import pandas as pd
import numpy as np
import matplotlib.pyplot as plt
import matplotlib.cm as cm
from sklearn.linear_model import Lasso
%matplotlib inline

path = os.getcwd() + "/7341_C2.txt"
data = pd.read_csv(path) # ,header=None
print(data.head())

cols = data.shape[1]
# convert from pandas dataframe to numpy matrices
X = np.matrix(data.iloc[:,1:cols].values)
y = np.array(data.iloc[:,0].values.ravel()) # first col is y label

# use map(float, ) to convert the string list to float list
X_names = list(map(float, data.columns.values[1:])) # X_names = np.array(list(data)[1:])
labels = list(set(y))

m = X.shape[0]
n = X.shape[1]

      Label  250   251   252   253   254        255        256        257 \
0       1     10     10     10     10     10    43.570310   31.487300  16.338870
1       1     10     10     10     10     10    6.282227   -4.985352  -6.870117
2       1     10     10     10     10     10   129.908200  123.578100  92.240230
3       1     10     10     10     10     10   105.136700   93.134770  60.101560
4       1     10     10     10     10     10   66.738280   68.101560  64.845700

      258    ...  2330  2331  2332  2333  2334  2335  2336  2337  2338  2339
0  10.359380  ...    10     10     10     10     10     10     10     10     10     10
1   2.167969  ...    10     10     10     10     10     10     10     10     10     10
2   50.460940  ...    10     10     10     10     10     10     10     10     10     10
3   18.257810  ...    10     10     10     10     10     10     10     10     10     10
4   58.904300  ...    10     10     10     10     10     10     10     10     10     10

[5 rows x 2091 columns]
```

Visualize Data

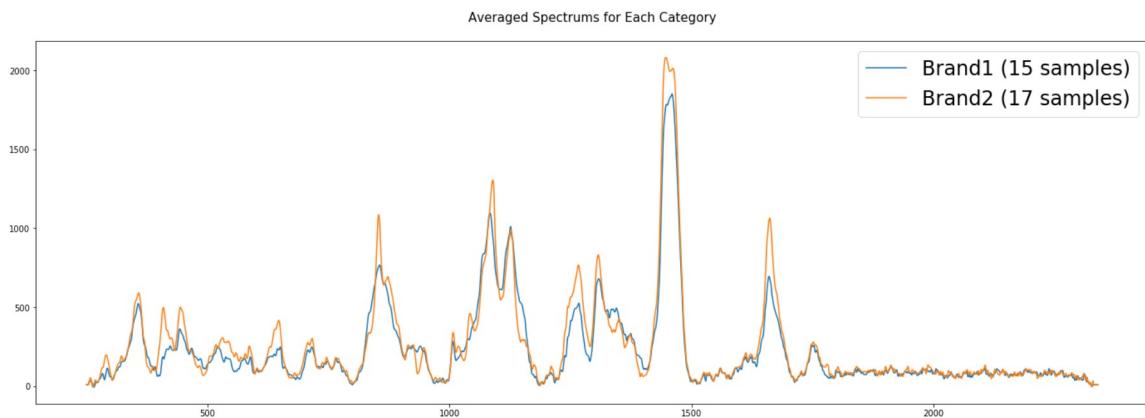
```
In [9]: import pylab, matplotlib
matplotlib.rcParams.update({'font.size': 18})

plt.figure(figsize = (24,8))

for c in labels:
    Xc = X[y == c]
    yc = y[y == c]
    plt.plot(X_names, np.mean(Xc, axis=0).tolist()[0], label= 'Brand' + str(c) +
' (' + str(len(yc)) + ' samples)')
    plt.legend()

plt.title(u'Averaged Spectrums for Each Category\n')
```

Out[9]: Text(0.5, 1.0, 'Averaged Spectrums for Each Category\n')



Plot the first two principal components in the 2D plane

```
In [11]: # %load plotComponents2D.py
import matplotlib.pyplot as plt
import numpy as np

def plotComponents2D(X, y, labels, use_markers = False, ax=None):

    if X.shape[1] < 2:
        print('ERROR: X MUST HAVE AT LEAST 2 FEATURES/COLUMNS! SKIPPING plotComponents2D().')
        return

    # Gray shades can be given as a string encoding a float in the 0-1 range
    colors = ['0.9', '0.0', '0.5', 'red', 'blue', 'black', 'orange', 'green', 'cyan', 'purple']
    markers = ['o', 's', '^', 'D', 'H', 'o', 's', '^', 'D', 'H', 'o', 's', '^', 'D', 'H']

    if (ax is None):
        fig, ax = plt.subplots()

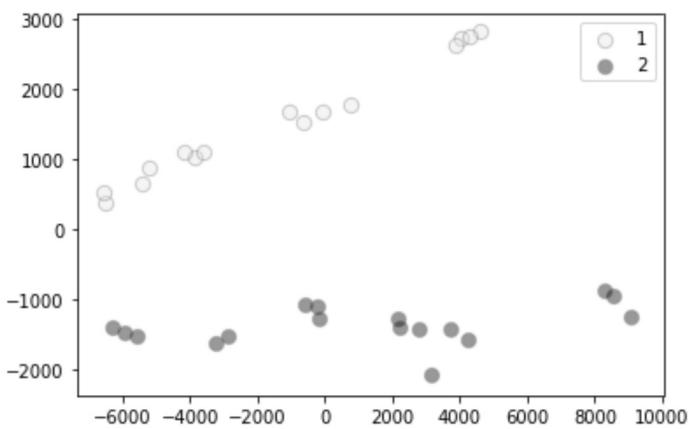
    i=0
    for label in labels:
        cluster = X[np.where(y == label)]
        # print(cluster.shape)

        if use_markers:
            ax.scatter([cluster[:,0]], [cluster[:,1]], s=40, marker=markers[i], facecolors='none', edgecolors=colors[i+3], label=str(label))
        else:
            ax.scatter([cluster[:,0]], [cluster[:,1]], s=70, facecolors=colors[i], label=str(label), edgecolors = 'gray', alpha = .4) # cmap='tab20'
        i=i+1

    ax.legend()
```

```
In [12]: from sklearn.decomposition import PCA

pca = PCA(n_components=2) # keep the first 2 components
pca.fit(X)
X_pca = pca.transform(X)
plotComponents2D(X_pca, y, labels)
```



The two classes have a wide dividing area. After CS, we want to keep this discriminative structure between the two classes.

Generating Sensing Matrix

```
In [51]: def getSensingMatrix(N, a = 0.2): # N = data set size
    # np.random.seed(10)

    OMEGA = np.zeros((N,N))
    pm = np.random.permutation(np.arange(0, N))
    for i in range(N):
        OMEGA[i, pm[i]] = 1

    PHI = OMEGA[:int(N*a)]

    return PHI, OMEGA
```

```
In [52]: getSensingMatrix(5, 0.2)
```

```
Out[52]: (array([[0., 0., 0., 1., 0.]]), array([[0., 0., 0., 1., 0.],
       [1., 0., 0., 0., 0.],
       [0., 0., 0., 0., 1.],
       [0., 0., 1., 0., 0.],
       [0., 1., 0., 0., 0.]]))
```

```
In [26]: # Another version of sensing /sampling. Without using PHI
```

```
def sampler(x, a = 0.2): # N = data set size, a - sampling ratio
    np.random.seed(10)
    N = len(x)
    pm = np.random.permutation(np.arange(1, N))
    r = pm[:int(N*a)]
    xs = x[r]
    return xs
```

```
In [305]: # the new version of sensing /sampling
```

```
def sensing(x, PHI): # x is a list
    assert (x.shape[0] == PHI.shape[1])
    return PHI @ x
```

```
In [308]: #x = np.random.randn(4)
# x, sampler(x, a = 0.5)

phi, omega = getSensingMatrix(10, 0.2)
x = np.random.randn(10)
x.shape, phi.shape, x, sensing(x, phi)
```

```
Out[308]: ((10,),
(2, 10),
array([-1.7639635 ,  0.84521648,  2.03049231, -0.55314797, -0.36917264,
       0.26074389, -1.12290167,  0.55530285, -0.51296706,  2.1098513 ],
array([0.84521648, 2.03049231]))
```

Choose Basis Matrix Ψ

Ψ is a unitary matrix .i.e. its conjugate transpose is also its inverse, or $\Psi^H = \Psi^{-1}$

For real matrices, unitary is the same as orthogonal .

Use 5 candidate basis matrices: IDM (Identity matrix), DCT (discrete cosine transform), DFT (which is widely used in medical imaging), HWT (Hadamard-Walsh transform, which is widely used in signal processing), UDM (uniformly distributed random orthogonal matrix)

IDM doesn't transform into other spaces. It is just for theoretical purpose.

DCT

The function `dctmtx()` generates the n-by-n discrete cosine transform (DCT) matrix.

$$DCT_{M,N} = \frac{1}{\sqrt{N}} \begin{bmatrix} 1 & 1 & \cdots & 1 \\ \sqrt{2}\cos\left(\frac{\pi}{2N}\right) & \sqrt{2}\cos\left(\frac{3\pi}{2N}\right) & \cdots & \sqrt{2}\cos\left(\frac{(2N-1)\pi}{2N}\right) \\ \vdots & \vdots & \ddots & \vdots \\ \sqrt{2}\cos\left(\frac{(M-1)\pi}{2N}\right) & \sqrt{2}\cos\left(\frac{3(M-1)\pi}{2N}\right) & \cdots & \sqrt{2}\cos\left(\frac{(2N-1)(M-1)\pi}{2N}\right) \end{bmatrix}$$

```
In [6]: import scipy

def dctmtx(m, n):

    mtx = np.zeros((m,n))
    N = n

    mtx[0, :] = 1 * np.sqrt(1/N)
    for i in range(1, m):
        for j in range(n):
            mtx[i, j] = np.cos(np.pi * i * (2*j+1) / (2*N)) * np.sqrt(2/N)

    plt.figure()
    plt.imshow(mtx, interpolation='nearest', cmap=cm.Greys_r)
    plt.axis('off')

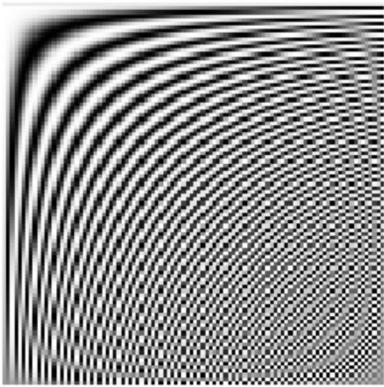
    print(np.fabs(scipy.linalg.det(mtx)))

    return mtx
```

Let us generate and visualize a 100x100 DCT matrix.

```
In [7]: dctmtx(100,100)
```

```
Out[7]: array([[ 0.1           ,  0.1           ,  0.1           , ...,  0.1           ,
   0.1           ,  0.1           ],
 [ 0.14140391,  0.14126436,  0.1409854 , ..., -0.1409854 ,
 -0.14126436, -0.14140391],
 [ 0.14135157,  0.14079372,  0.13968022, ...,  0.13968022,
  0.14079372,  0.14135157],
 ...,
 [ 0.00666186, -0.01992644,  0.03301416, ..., -0.03301416,
  0.01992644, -0.00666186],
 [ 0.00444215, -0.01330893,  0.02212317, ...,  0.02212317,
 -0.01330893,  0.00444215],
 [ 0.00222135, -0.00666186,  0.01109579, ..., -0.01109579,
  0.00666186, -0.00222135]])
```



DFT

The function `dftmtx()` generate the n-by-n discrete fourier transform (DFT) matrix

$$DFT_{N,N} = \left(\frac{\omega^{jk}}{\sqrt{N}} \right)_{j,k=0,\dots,N-1} = \frac{1}{\sqrt{N}} \begin{bmatrix} \omega^0 & \omega^0 & \omega^0 & \cdots & \omega^0 \\ \omega^0 & \omega^1 & \omega^2 & \cdots & \omega^{N-1} \\ \omega^0 & \omega^2 & \omega^4 & \cdots & \omega^{2(N-1)} \\ \omega^0 & \omega^3 & \omega^6 & \cdots & \omega^{3(N-1)} \\ \vdots & \vdots & \vdots & \ddots & \vdots \\ \omega^0 & \omega^{N-1} & \omega^{2(N-1)} & \cdots & \omega^{(N-1)(N-1)} \end{bmatrix}$$

where

$$\omega = e^{-2\pi i/N} = \cos(-2\pi/N) + i\sin(-2\pi/N)$$

```
In [8]: import numpy as np
import scipy

def dftmtx(N):
    i, j = np.meshgrid(np.arange(N), np.arange(N))
    w = np.exp(-2 * np.pi * 1j / N)
    mtx = np.power(w, i * j) / np.sqrt(N)

    plt.figure(figsize=(12, 6))
    # plt.title("abs | phase")
    plt.subplot(1, 2, 1)
    plt.imshow(abs(mtx), interpolation='nearest', cmap=cm.Greys_r)
    plt.axis('off')
    plt.subplot(1, 2, 2)
    plt.imshow(np.angle(mtx), interpolation='nearest', cmap=cm.Greys_r)
    plt.axis('off')

    print(scipy.linalg.det(mtx))

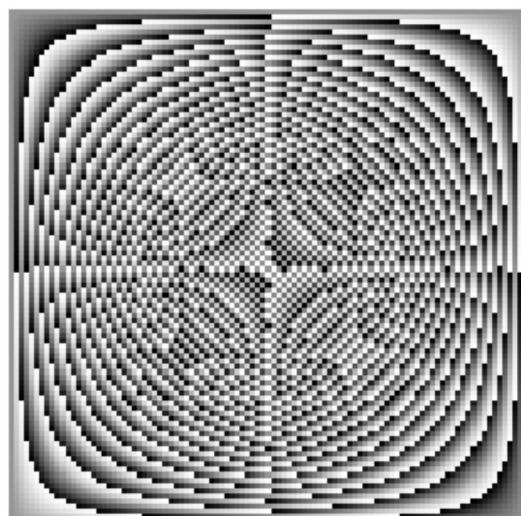
    return mtx
```

Let us generate a 100x100 DFT matrix, and visualize the amplitude (abs) and phase (angle) components.

```
In [9]: dftmtx(100)

(7.477907182362742e-13+0.9999999999993605j)
```

```
Out[9]: array([[0.1          +0.j        , 0.1          +0.j        ,
   0.1          +0.j        , ..., 0.1          +0.j        ,
   0.1          +0.j        , 0.1          +0.j        ],
  [0.1          +0.j        , 0.09980267-0.00627905j,
   0.09921147-0.01253332j, ..., 0.09822873+0.01873813j,
   0.09921147+0.01253332j, 0.09980267+0.00627905j],
  [0.1          +0.j        , 0.09921147-0.01253332j,
   0.09685832-0.02486899j, ..., 0.09297765+0.03681246j,
   0.09685832+0.02486899j, 0.09921147+0.01253332j],
  ...,
  [0.1          +0.j        , 0.09822873+0.01873813j,
   0.09297765+0.03681246j, ..., 0.08443279-0.05358268j,
   0.09297765-0.03681246j, 0.09822873-0.01873813j],
  [0.1          +0.j        , 0.09921147+0.01253332j,
   0.09685832+0.02486899j, ..., 0.09297765-0.03681246j,
   0.09685832-0.02486899j, 0.09921147-0.01253332j],
  [0.1          +0.j        , 0.09980267+0.00627905j,
   0.09921147+0.01253332j, ..., 0.09822873-0.01873813j,
   0.09921147-0.01253332j, 0.09980267-0.00627905j]])
```



HWT

The Hadamard-Walsh transform (HWT) matrix requires the dimension to be 2^N .

In function `hwtmtx()`, we will round the input parameter to its next 2^N value.

$$HWT_{(2^N)} = \frac{1}{\sqrt{N/2}} \begin{bmatrix} 1 & 1 & 1 & 1 & \cdots & 1 & 1 \\ 1 & -1 & 1 & -1 & \cdots & 1 & -1 \\ 1 & 1 & 1 & 1 & \cdots & 1 & 1 \\ 1 & -1 & 1 & -1 & \cdots & 1 & -1 \\ \vdots & \vdots & \vdots & \vdots & \ddots & \vdots & \vdots \\ 1 & 1 & 1 & 1 & \cdots & 1 & 1 \\ 1 & -1 & 1 & -1 & \cdots & 1 & -1 \end{bmatrix}$$

```
In [14]: from scipy.linalg import hadamard
import math

def hwtmtx(N):

    NN = 2** (N-1).bit_length() # make sure it is a 2**n value
    print('Expanded to ', NN, ', Divide by', 2** (math.log(NN,2)/2))

    mtx = hadamard(NN) / (2** (math.log(NN,2)/2))

    plt.figure()
    plt.imshow(mtx, interpolation='nearest', cmap=cm.Greys_r)
    plt.axis('off')

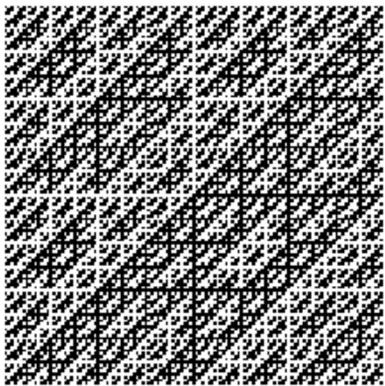
    # Use slogdet when mtx is big, e.g. > 2000
    logdet = np.linalg.slogdet(mtx)
    det = logdet[0] * np.exp(logdet[1])
    print('det =', det)

    return mtx, NN
```

```
In [15]: hwtmtx(100)
```

```
Expanded to 128 , Divide by 11.313708498984761
det = 0.99999999999999938
```

```
Out[15]: (array([[ 0.08838835,  0.08838835,  0.08838835, ...,  0.08838835,
   0.08838835,  0.08838835],
  [ 0.08838835, -0.08838835,  0.08838835, ..., -0.08838835,
   0.08838835, -0.08838835],
  [ 0.08838835,  0.08838835, -0.08838835, ...,  0.08838835,
  -0.08838835, -0.08838835],
  ...,
  [ 0.08838835, -0.08838835,  0.08838835, ...,  0.08838835,
  -0.08838835,  0.08838835],
  [ 0.08838835,  0.08838835, -0.08838835, ..., -0.08838835,
   0.08838835,  0.08838835],
  [ 0.08838835, -0.08838835, -0.08838835, ...,  0.08838835,
   0.08838835, -0.08838835]]), 128)
```



In the above example, hwtmtx(100) generates a 128x128 matrix. $128 = 2^7$

UDM (uniformly distributed random orthogonal matrix)

Function `rvsmtx()` generates a uniformly distributed random orthogonal matrix.

```
In [20]: from scipy.stats import ortho_group

def rvsmtx(N):
    mtx = ortho_group.rvs(N) # uniformly distributed random orthogonal matrix

    plt.figure()
    plt.imshow(mtx, interpolation='nearest', cmap=cm.Greys_r)
    plt.axis('off')

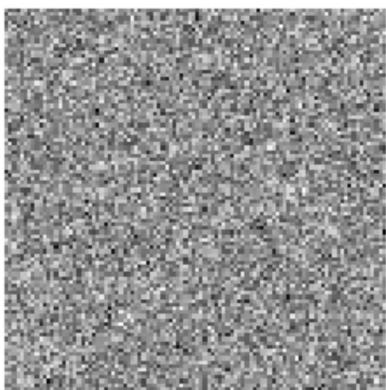
    print(np.fabs(scipy.linalg.det(mtx)))

    return mtx
```

In [21]: `rvsmtx(100)`

```
1.0000000000000044
```

Out[21]: `array([[0.02662416, -0.14342694, -0.10566114, ..., -0.08841288,
 0.16754891, 0.02157665],
 [-0.17249476, 0.14768207, 0.06745075, ..., 0.11603294,
 -0.00545798, 0.02949385],
 [-0.0241674 , -0.09139529, -0.06255712, ..., 0.05553968,
 0.04324998, -0.00282775],
 ...,
 [-0.00478052, 0.05895622, -0.06395684, ..., 0.27288879,
 0.03907859, 0.02712428],
 [-0.10277921, -0.10599319, 0.0207901 , ..., 0.03194886,
 -0.09970206, -0.00548009],
 [0.04876354, 0.06733024, -0.08274201, ..., 0.11713795,
 0.08942622, 0.05651709]])`



Generate 5 candidate Ψ , and save to PSIS.pkl file.

Next time we can just deserialize the pkl file to load back.

In []: `PSIS=[
 np.identity(n), # the identity matrix, just for comparison
 dctmtx(n,n),
 dftmtx(n),
 hwtmtx(n)[0],
 rvsmtx(n)
]

import pickle

filehandler = open("PSIS.pkl", "wb")
pickle.dump(PSIS,filehandler)
filehandler.close()`

Load PSIs back from pickle

In [22]: `# Load PSIs back from pickle

import pickle

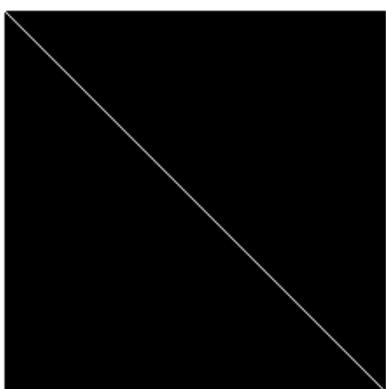
file = open("PSIS.pkl",'rb')
PSIS = pickle.load(file)
file.close()`

Visualize the Ψ candidates

```
In [25]: import matplotlib.pyplot as plt
import matplotlib.cm as cm
%matplotlib inline

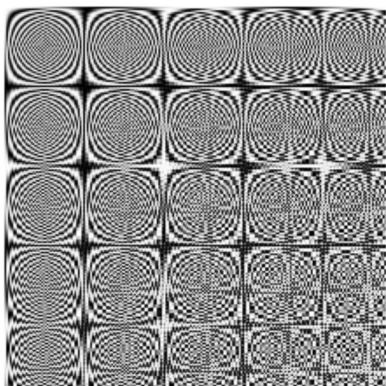
plt.figure()
plt.imshow(PSIS[0], interpolation='nearest', cmap=cm.Greys_r)
plt.axis('off')
print('IDM')
```

IDM



```
In [27]: plt.figure()
plt.imshow(PSIS[1], interpolation='nearest', cmap=cm.Greys_r)
plt.axis('off')
print('DCT')
```

DCT

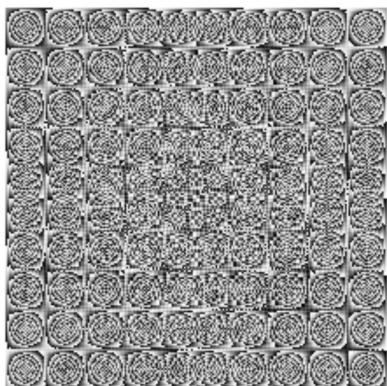
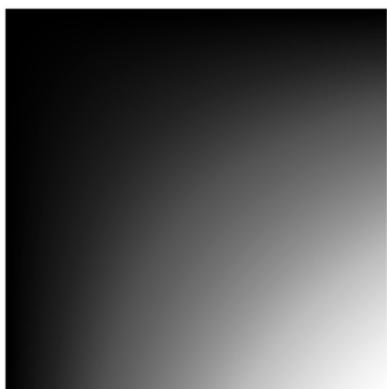


```
In [26]: import numpy as np
```

```
plt.figure()
plt.imshow(abs(PSIS[2]), interpolation='nearest', cmap=cm.Greys_r)
plt.axis('off')

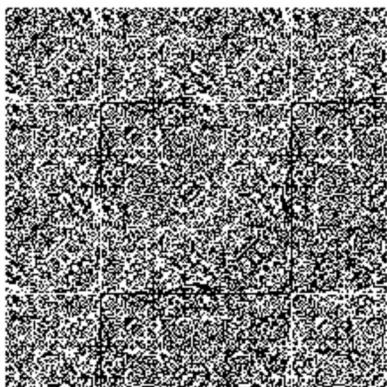
plt.figure()
plt.imshow(np.angle(PSIS[2]), interpolation='nearest', cmap=cm.Greys_r)
plt.axis('off')
print('DFT')
```

DFT



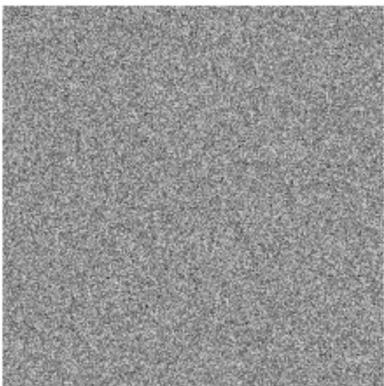
```
In [32]: plt.figure()
plt.imshow(PSIS[4], interpolation='nearest', cmap=cm.Greys_r)
plt.axis('off')
print('HWT')
```

HWT



```
In [33]: plt.figure()
plt.imshow(PSIS[5], interpolation='nearest', cmap=cm.Greys_r)
plt.axis('off')
print('UDM')
```

UDM



Coherence Between Two Bases

$$\mu(A, B) = \sqrt{n} \max_{0 \leq i, j < n} (a_i^T b_j)$$

a and b are column vectors in A and B

$$0 \leq \mu(A, B) \leq \sqrt{n}$$

We are interested in bases pairs with small coherence and in that case we say the two bases are incoherent. In the case of dictionaries for sparse representation, the incoherence is important because a small μ means small correlation between the two bases, thus a dictionary composed by such a pair of bases has a “richer vocabulary” relative to a dictionary with a larger μ .

CS requires a small μ between Ψ and Ω . Ω is a scrambled identity matrix. The sensing matrix Φ is a subset off Ω (take the first few rows).

Theoretical computation results:

$$\begin{aligned}\mu(\Omega, I) &= \sqrt{n} \\ \mu(\Omega, C) &= \sqrt{2} \\ \mu(\Omega, F) &= 1 \\ \mu(\Omega, H) &= 1\end{aligned}$$

$$I = IDM, C = DCT, F = DFT, H = HWT$$

Function `mutual_coherence()` is defined to calculate μ .

We will check whether the experiment result conforms to the theoretical result.

```
In [49]: from tqdm import tqdm

def mutual_coherence(A,B):
    assert (A.shape == B.shape)
    assert (A.shape[0] == A.shape[1])
    n = A.shape[0]
    p = 0

    with tqdm(total= n*n) as pbar:
        for i in range(n):
            for j in range(n):
                a = A[:,i].T
                b = B[:,j]
                tmp = a@b
                # print(tmp)
                if (p < tmp):
                    p = tmp
                pbar.update(1)
    return math.sqrt(n)*p
```

```
In [319]: psi_names = ['Identity Matrix', 'DCT', 'DFT', 'DWT', 'Hadamard-Walsh Matrix', 'Uniformly Distributed Random Orthogonal Matrix']

for idx, PSI in enumerate(PSIS):
    assert (PSI.shape[0] == PSI.shape[1])

    # for some transformations, PSI dimension may differ. e.g. HWT requires 2**n and DWT requires even number
    n = PSI.shape[0]
    _, OMEGA = getSensingMatrix(n)

    print(psi_names[idx], ":", mutual_coherence(PSI, OMEGA))
```

100% |██████████| 43
68100/4368100 [03:48<00:00, 19152.96it/s]

Identity Matrix : 45.71651780264984

100% |██████████| 43
68100/4368100 [03:48<00:00, 19156.82it/s]

DCT : 1.4142135623730951

100% |██████████| 43
68100/4368100 [03:10<00:00, 22978.39it/s]

DFT : (1.00000000006146+7.877968100069424e-13j)

100% |██████████| 1677
7216/16777216 [16:58<00:00, 16469.66it/s]

Hadamard-Walsh Matrix : 1.0

100% |██████████| 43
68100/4368100 [03:48<00:00, 19138.18it/s]

Uniformly Distributed Random Orthogonal Matrix : 5.293445504102594

The result perfectly matches the theoretic calculation.

IDM reaches the upper bound of μ , i.e. \sqrt{n}

DCT: $\sqrt{2}$

DFT: 1

Hadamard-Walsh Matrix: 1

Sparsity analysis

$$z = \Psi^H x$$

We will transform the original x into the latent space defined by Ψ .

CS requires to choose the Ψ that results in the highest sparsity.

```
In [601]: x = X[0,:]
x.shape

import pylab, matplotlib

plt.figure(figsize=(24,48))
rows = len(PSIS)
matplotlib.rcParams.update({'font.size': 18})

for idx, PSI in enumerate(PSIS):

    xx = np.copy(x)

    # pad x with zero if necessary
    if (x.shape[1] < PSI.shape[0]):
        xx = np.zeros((x.shape[0], PSI.shape[0]))
        xx[:x.shape[0], :x.shape[1]] = x

    # print(xx.shape, PSI.shape)

    PSI_H = PSI.conj().T
    PSI_INV = np.linalg.pinv(PSI)
    #print(PSI_H)
    #print(PSI_INV)
    # theoretically, PSI_H == PSI_INV
    z = (xx @ PSI_H).ravel().tolist()
    # print(z)
    MAX = abs(max(z, key=abs))
    print(MAX)
    thresholds = np.array(range(100)) / 5000

    rs = []
    for threshold in thresholds:
        rs.append((np.abs(np.array(z)) <= threshold * MAX).sum() / len(z))

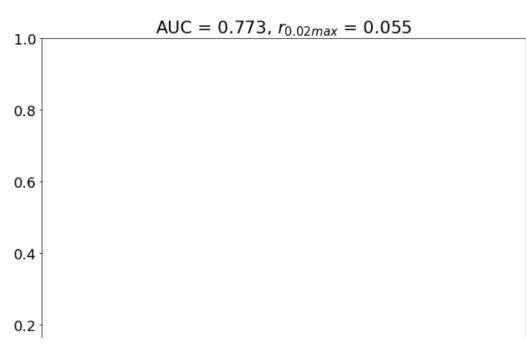
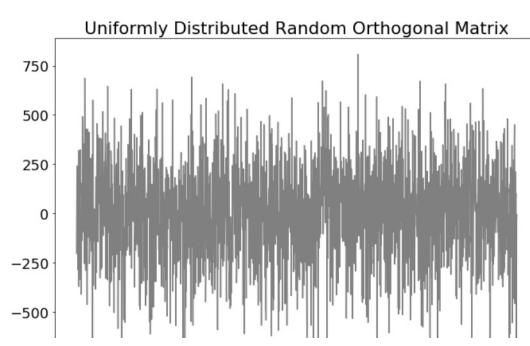
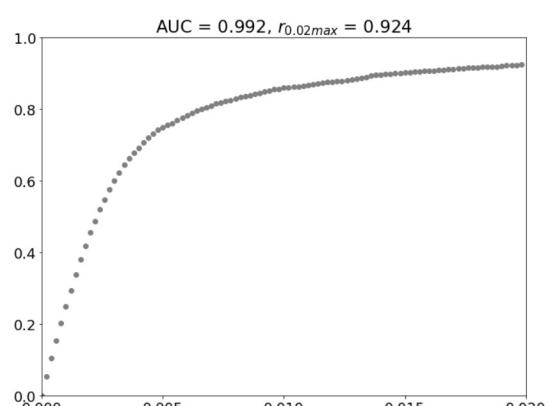
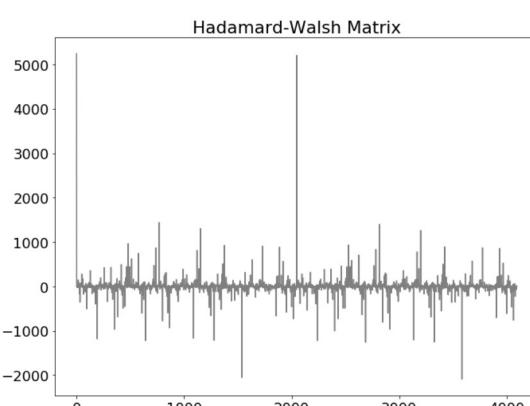
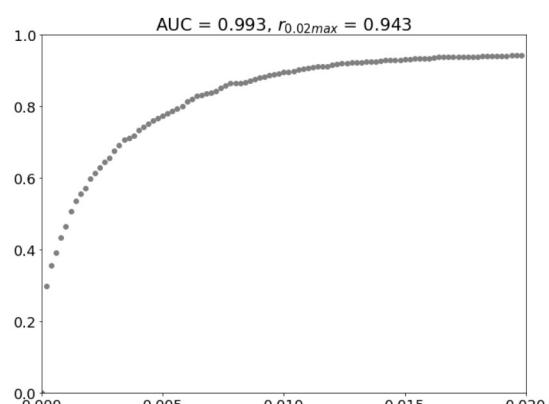
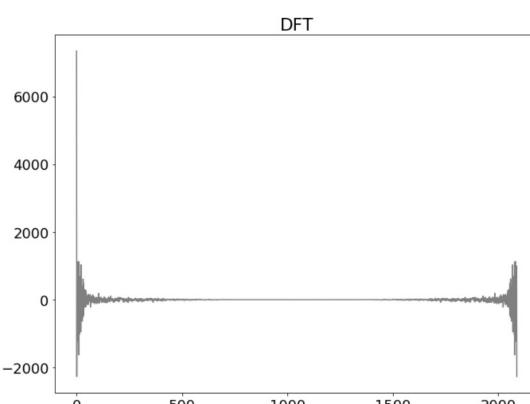
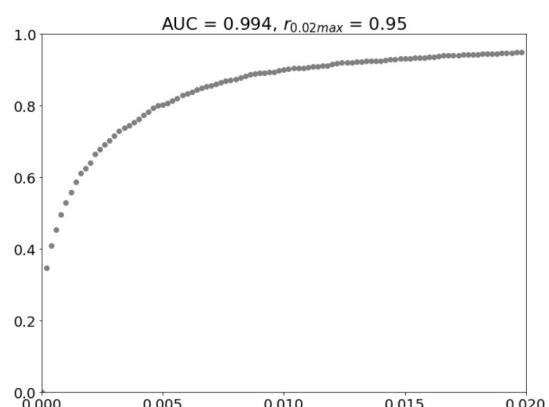
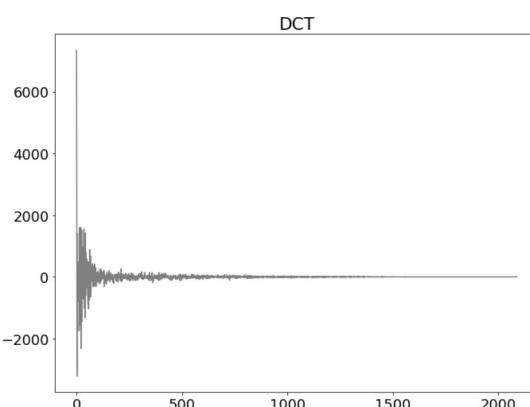
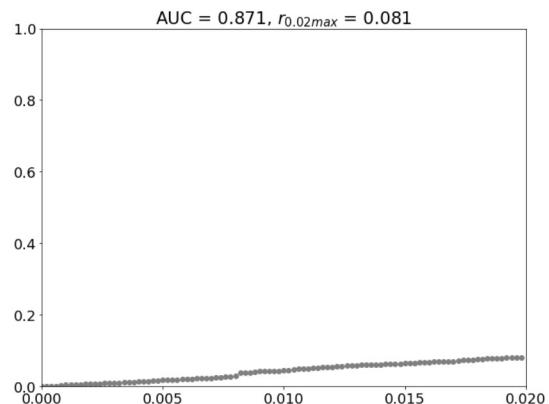
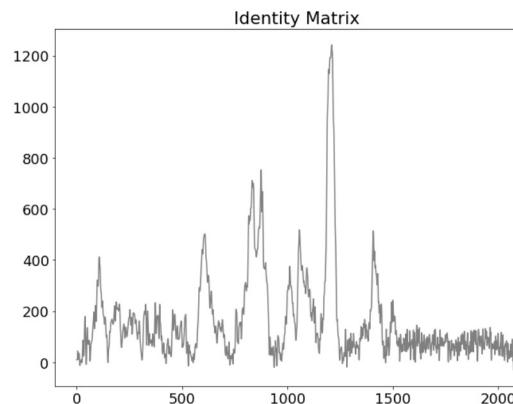
    auc = 0
    for i in range(1000):
        auc += (np.abs(np.array(z)) <= (i+1)/1000 * MAX).sum() / len(z)
    auc = auc/1000

    r = (np.abs(np.array(z)) <= 0.02 * MAX).sum() / len(z) # use 0.02 MAX ABS
    as threshold

    plt.subplot(rows,2,2*idx+1)
    plt.plot(z, color='gray')
    plt.title(psi_names[idx])
    # plt.axis('off')

    plt.subplot(rows,2,2*idx+2)
    plt.scatter(thresholds, rs, color='gray')
    plt.title('AUC = ' + str(round(auc,3)) + ', $r_{0.02max} = ' + str(round(r,3)))
    pylab.xlim([0,0.02])
    pylab.ylim([0,1.0])
    plt.xticks(np.arange(0.0, 0.021, 0.005))
```

1243.019
7338.632439548879
7338.632439548878
5242.136258902184
845.6028616661284



Best AUC: DFT DCT
Best μ : DFT HWT

Finally, Choose DFT

Under DFT, at the 0.02 max abs threshold, non-sparsity degree $d \approx 1 - 0.95 = 0.05$. According to the "four-to-one" practical / empirical rule: for exact reconstruction, one needs about four incoherent measurements per unknown nonzero term

$k \geq 4d$ (d is percentage of non-zeros in the signal, 1 - degree of sparsity)
 $k \geq 0.2$, i.e. 0.2 is the lower bound for effective CS.

Sensing Matrix for sub-Nyquist sampling

batch process by vectorization

```
In [348]: k = 0.2
PHI, OMEGA = getSensingMatrix(X.shape[1], k)

PHI.shape, OMEGA.shape
```

```
Out[348]: ((418, 2090), (2090, 2090))
```

```
In [349]: Xs = PHI @ X.T
Xs.shape
```

```
Out[349]: (418, 46)
```

single-sample processing

```
In [55]: import numpy as np

def sampler(x, k = 0.2): # N = data set size, a - sampling ratio

    N = len(x)

    # when k > 1, return x directly
    if (k > 1):
        return x, list(range(N))

    pm = np.random.permutation(np.arange(1, N))
    idx = pm[:int(N*k)]
    xs = np.array(x)[idx]

    return xs, idx
```

```
In [387]: x = X[0,:].ravel().tolist()[0] # get the first sample
xs, idx = sampler(x, k = 0.2)
xs.shape, idx
```

```
Out[387]: ((418,),  
 array([1105, 1875, 539, 141, 1973, 1071, 1048, 577, 783, 5, 1753,  
 1687, 1133, 1950, 1443, 808, 1634, 2085, 928, 1879, 1683, 1736,  
 1448, 1888, 2038, 1178, 1053, 460, 2042, 1150, 992, 462, 919,  
 405, 1198, 1589, 862, 1721, 768, 731, 1580, 1974, 107, 644,  
 1077, 274, 28, 1792, 1242, 263, 1026, 1598, 1394, 293, 1819,  
 2004, 1136, 259, 642, 1060, 629, 612, 1384, 2039, 9, 192,  
 1309, 1926, 203, 1931, 1465, 1109, 1392, 1829, 334, 1254, 1887,  
 1399, 907, 1708, 195, 1203, 1821, 483, 1230, 247, 340, 690,  
 181, 1986, 261, 27, 666, 354, 1115, 1741, 682, 601, 296,  
 1776, 39, 1604, 1345, 1939, 89, 1415, 1471, 110, 1353, 1566,  
 498, 574, 1765, 98, 760, 1251, 1014, 212, 1482, 1218, 1913,  
 1354, 2028, 670, 278, 14, 1565, 356, 1696, 8, 1045, 1749,  
 774, 267, 1331, 1219, 943, 708, 871, 991, 1395, 667, 1393,  
 464, 38, 1861, 1369, 1293, 843, 1490, 398, 132, 1632, 924,  
 297, 438, 57, 16, 540, 1593, 1344, 338, 1912, 73, 1599,  
 108, 791, 439, 359, 1483, 990, 1356, 1576, 578, 741, 1185,  
 517, 1426, 139, 1489, 1383, 1248, 295, 1911, 446, 1874, 163,  
 993, 1363, 222, 1271, 428, 2009, 81, 710, 1481, 766, 2082,  
 199, 1241, 288, 1863, 727, 887, 200, 53, 1575, 117, 223,  
 1153, 1275, 2019, 492, 1664, 789, 1731, 207, 273, 479, 1716,  
 1257, 635, 1584, 1449, 1639, 482, 1247, 1166, 1480, 1617, 1520,  
 672, 976, 1928, 905, 1364, 1311, 893, 812, 1917, 1260, 55,  
 883, 1336, 1312, 1085, 179, 1339, 1704, 617, 1027, 1952, 845,  
 586, 360, 282, 1327, 941, 1552, 37, 499, 1050, 2065, 1463,  
 129, 1979, 2035, 140, 759, 451, 720, 1243, 643, 886, 1496,  
 35, 916, 949, 1387, 403, 374, 1236, 1341, 1146, 1597, 765,  
 1838, 1557, 1603, 1088, 1832, 1036, 1194, 1944, 501, 1120, 364,  
 1648, 1854, 610, 685, 175, 563, 1382, 535, 1497, 614, 312,  
 1375, 860, 36, 399, 333, 1209, 101, 51, 1815, 281, 728,  
 477, 246, 82, 1823, 951, 372, 1074, 1684, 1529, 1372, 1608,  
 528, 299, 1102, 1359, 944, 1546, 2023, 30, 1763, 1280, 1714,  
 64, 1942, 136, 1017, 1197, 1056, 1335, 835, 343, 1184, 1833,  
 1154, 2048, 434, 1717, 1798, 233, 300, 1170, 1880, 622, 1945,  
 1622, 1836, 604, 1706, 549, 541, 1447, 1641, 668, 1761, 1637,  
 693, 1222, 1454, 1978, 1021, 347, 230, 750, 1234, 1029, 1971,  
 1065, 61, 821, 600, 489, 1627, 816, 1545, 189, 1245, 2079,  
 313, 840, 1672, 465, 1610, 769, 1317, 275, 873, 2016, 1667,  
 1445, 1286, 271, 1516, 879, 1840, 1884, 1573, 1340, 908, 1411]))
```

From the original 2090-point signal, only 418 points are sampled.

Measuring Matrix $A = \Phi\Psi$

```
In [361]: # Get the PSI for DFT
PSI = PSIS[1]
```

```
In [366]: A = PHI @ PSI
A.shape, PHI.shape, PSI.shape, Xs.shape
```

```
Out[366]: ((418, 2090), (418, 2090), (2090, 2090), (418, 46))
```

can also use the following to get the measuring matrix A

```
In [3]: import cv2
```

```
def csmtx(N, idx, t = 'dct'):

    K = len(idx)
    a = np.zeros((1, N))

    # Suppose the first number in the permutation is 42. Then choose the 42th
    # elements in a matrix full of zeros and change it to 1. Apply a discrete
    # cosine transformation to this matrix, and add the result to other matrix
    # Adelta as a new row.
    A = np.zeros((K, N))
    for i, j in enumerate(idx):
        a[0,j] = 1
        if t == 'dft':
            A[i, :] = cv2.dft(a)
        else:
            A[i, :] = cv2.dct(a)
        a[0,j] = 0

    ##### ALTERNATIVE IMPLEMENTATION #####
    # phi = np.zeros((K, N)) # Sampling matrix, kxN
    # for i, j in enumerate(r):
    #     phi[i,j] = 1
    # psi = dctmtx(N,N) # memory costly
    # A = phi*psi

    return A
```

```
In [391]: A = csmtx(len(x), idx, 'dft')
A.shape
```

```
Out[391]: (418, 2090)
```

L1 Minimization & Reconstruction

Let $A = \Phi\Psi$,

$$x_s = Az$$

$$A = [a_1, a_2, \dots, a_n]$$

Each a is column vector / atom / dictionary entry.

Because A has more columns than rows, A is overcomplete. That means A has excessive dictionary entries (non-repetitive), and can express data with high sparsity.

A is also full row rank, i.e. $\text{rank}(A) = r$ (row number)

The equations specified by $x_s = Az$ has many (not unique) solutions. This is an underdetermined linear system that has more unknowns than equations.

The solutions can be represented as $z = A^T(AA^T)^{-1}x_s + V[r :, :]v$, $A = USV^T$, v is a freely chosen $(n-r) \times 1$ vector.

We can find the sparsist solution by minimize L1-norm or $\|z\|_1$. We don't use L0-norm because minimizing it is NP hard.

```
objective: minimize ||z||
subject to: x_s = Az
```

```
In [49]: from sklearn.linear_model import Lasso
import cv2
import matplotlib

def lasso_cs(A, xs, L1, t = 'dct', silent = 'True'):
    lasso = Lasso(alpha=L1)
    lasso.fit(A, xs)
    # print ('non-zero coefs: ', np.count_nonzero(lasso.coef_))

    if t=='dft':
        x_r = cv2.idft(lasso.coef_)
    else:
        x_r = cv2.idct(lasso.coef_)

    if (silent == False):
        matplotlib.rcParams.update({'font.size': 20})
        fig, ax = plt.subplots(nrows=1, ncols=2, figsize=(32, 6))
        ax[0].plot(lasso.coef_, color = 'gray')
        #ax[0].title.set_text('z')

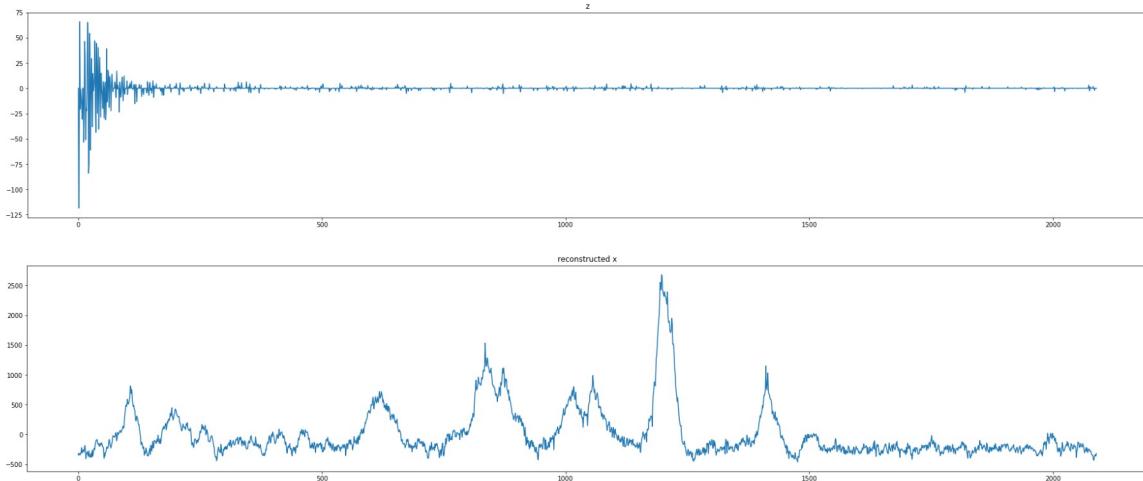
        ax[1].plot(x_r, color = 'gray')
        #ax[1].title.set_text('reconstructed x')

        fig.tight_layout()
        plt.show()

    return x_r, lasso.coef_
```

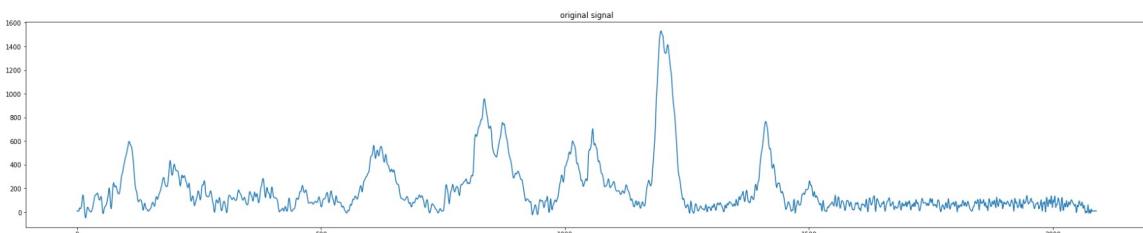
```
In [408]: xr, z = lasso_cs(A, xs, 0.005, 'dft', silent = False)
```

non-zero coefs: 621



```
In [410]: plt.figure(figsize=(32, 6))
plt.plot(x)
plt.title('original signal')
```

```
Out[410]: Text(0.5, 1.0, 'original signal')
```



Let us put all the above functions together

```
In [56]: def one_sample_CS(x, k = 0.2, L1 = 0.002, t='dft', silent = True):  
    xs, idx = sampler(x, k)  
    A = csmtx(len(x), idx, t)  
  
    xr, z = lasso_CS(A, xs, L1 = L1, t = t, silent = silent)  
  
    return xr, z
```

```
In [61]: x = X[0,:].ravel().tolist()[0]

PSI_H = dftmtx(len(x)).conj().T
z = (x @ PSI_H).ravel().tolist()

### Draw the original x ###
matplotlib.rcParams.update({'font.size': 20})
fig, ax = plt.subplots(nrows=1, ncols=2, figsize=(32, 6))
ax[0].plot(z, color = 'gray')
#ax[0].title.set_text('z')

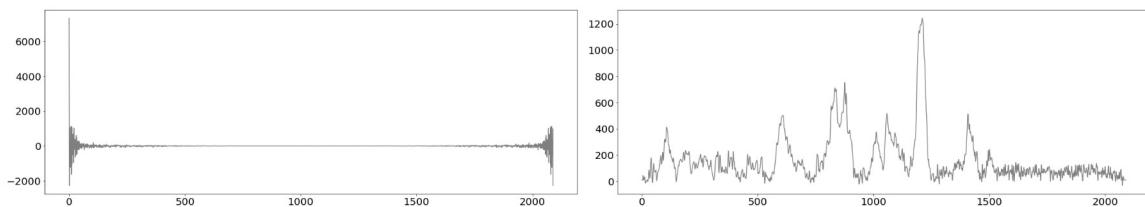
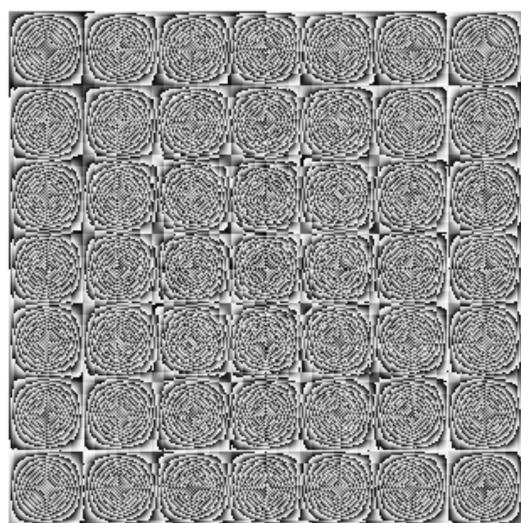
ax[1].plot(x, color = 'gray')
#ax[1].title.set_text('reconstructed x')

fig.tight_layout()
plt.show()

for k in [0.1, 0.2, 0.3, 0.4, 1.1]:
    for L1 in [0, 0.001, 0.01, 0.1, 1, 10]:
        print('===== k =', k, ', L1 =', L1, ' =====')
        one_sample_CS(x, k = k, L1 = L1, t='dft', silent = False)
```

```
(-1.0000000335595836-1.2473266863821664e-10j)
```

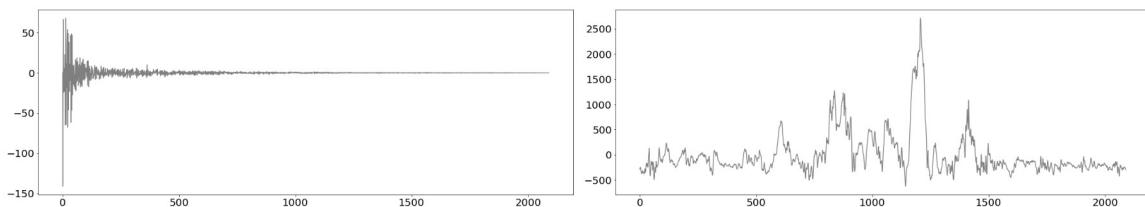
```
C:\Users\eleve\Anaconda3\lib\site-packages\numpy\core\numeric.py:538: ComplexWarning: Casting complex values to real discards the imaginary part
    return array(a, dtype, copy=False, order=order)
```



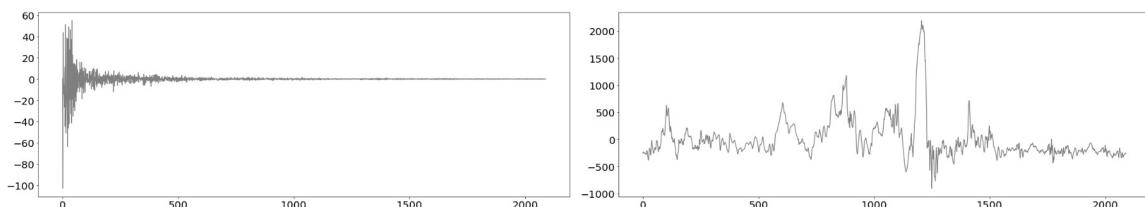
```
===== k = 0.1 , L1 = 0 =====
```

```
C:\Users\eleve\Anaconda3\lib\site-packages\ipykernel_launcher.py:7: UserWarning: With alpha=0, this algorithm does not converge well. You are advised to use the LinearRegression estimator
```

```
    import sys
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: UserWarning: Coordinate descent with no regularization may lead to unexpected results and is discouraged.
    positive)
```

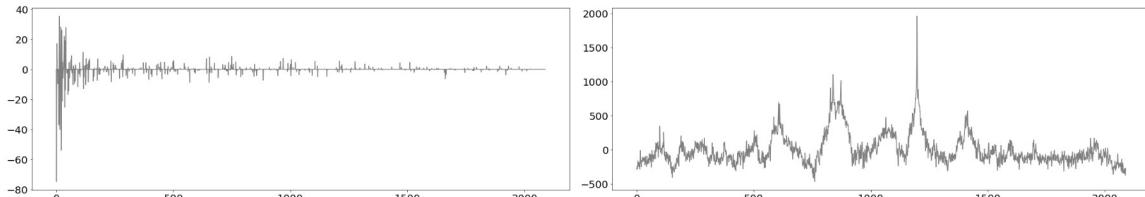


```
===== k = 0.1 , L1 = 0.001 =====
```



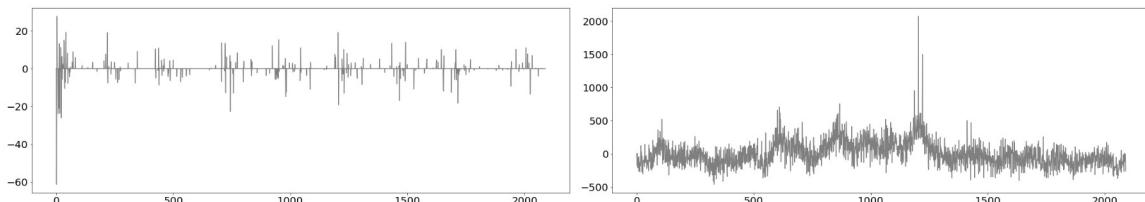
```
===== k = 0.1 , L1 = 0.01 =====
```

```
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 948.060011874864, tolerance: 4
41.08025135244895
    positive)
```

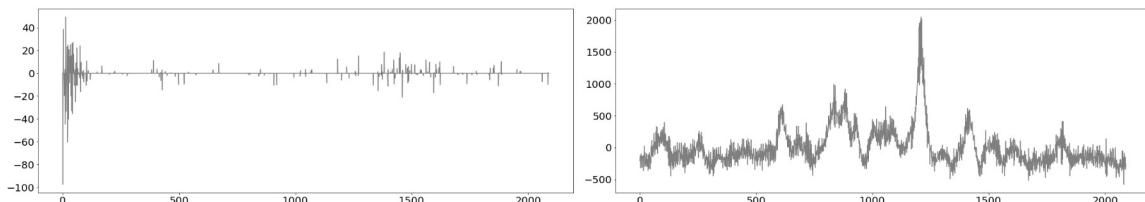


===== k = 0.1 , L1 = 0.1 =====

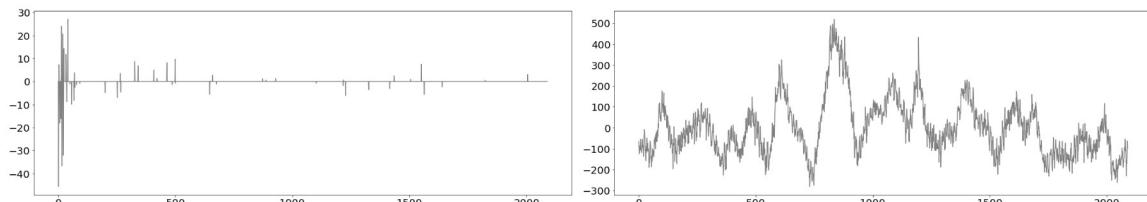
```
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 905.6630159714474, tolerance: 423.1507963585478  
    positive)
```



===== k = 0.1 , L1 = 1 =====



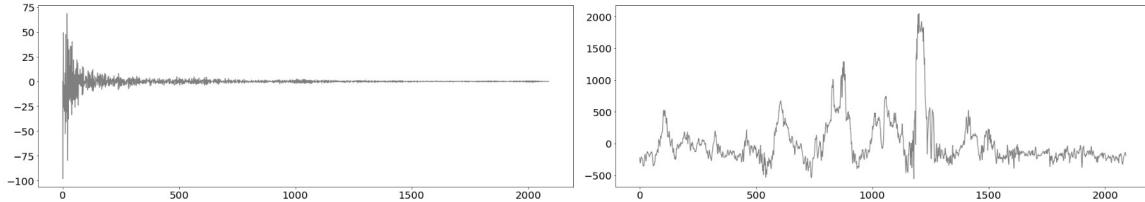
===== k = 0.1 , L1 = 10 =====



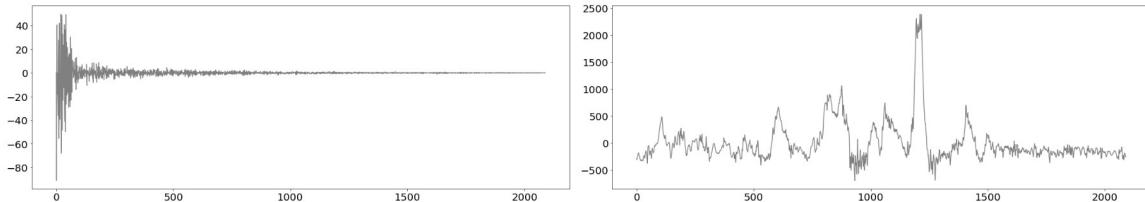
===== k = 0.2 , L1 = 0 =====

```
C:\Users\eleve\Anaconda3\lib\site-packages\ipykernel_launcher.py:7: UserWarning: With alpha=0, this algorithm does not converge well. You are advised to use the LinearRegression estimator
```

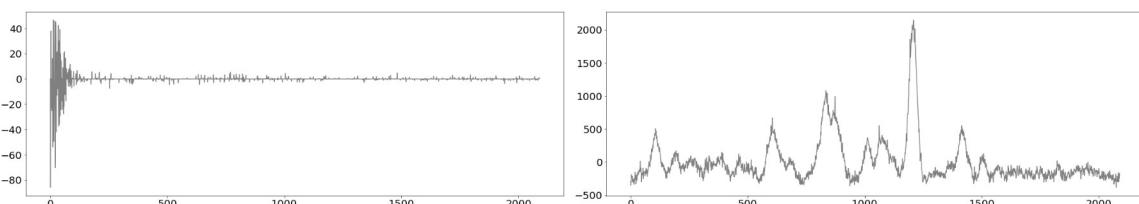
```
    import sys  
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: UserWarning: Coordinate descent with no regularization may lead to unexpected results and is discouraged.  
    positive)
```



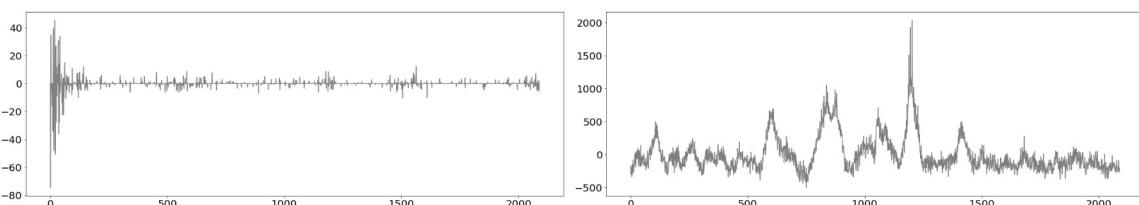
===== k = 0.2 , L1 = 0.001 =====



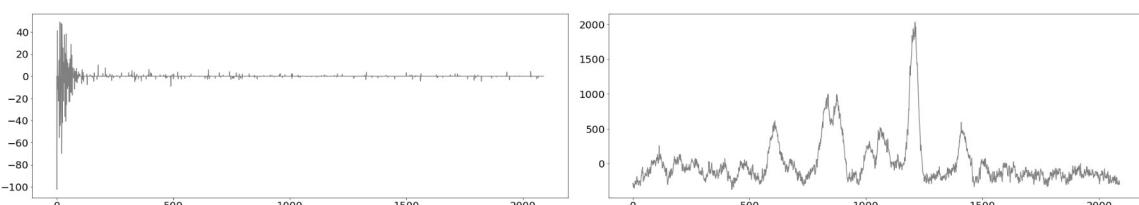
===== k = 0.2 , L1 = 0.01 =====



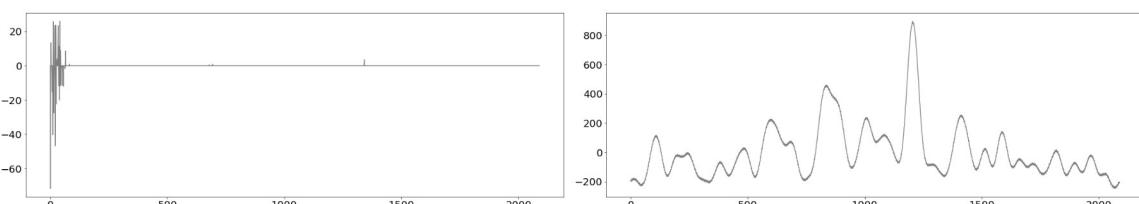
===== k = 0.2 , L1 = 0.1 =====



===== k = 0.2 , L1 = 1 =====

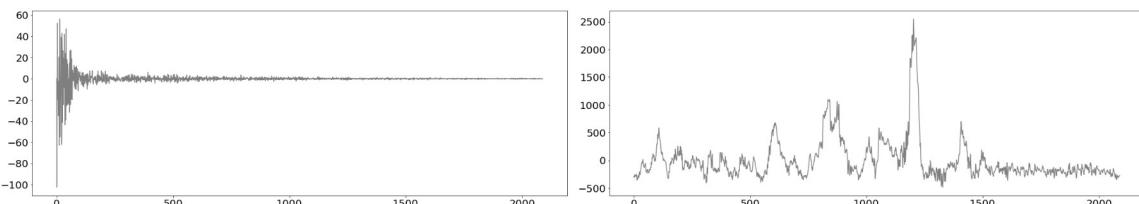


===== k = 0.2 , L1 = 10 =====

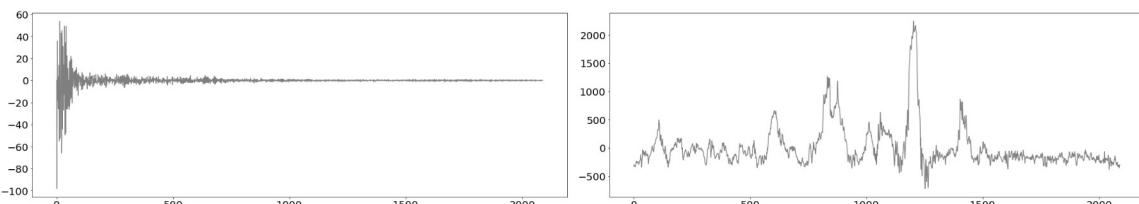


===== k = 0.3 , L1 = 0 =====

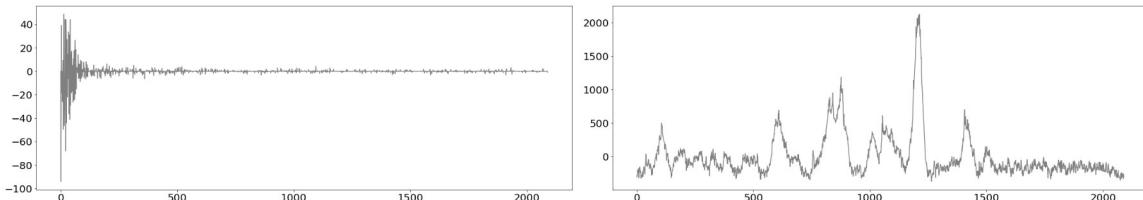
```
C:\Users\eleve\Anaconda3\lib\site-packages\ipykernel_launcher.py:7: UserWarning: With alpha=0, this algorithm does not converge well. You are advised to use the LinearRegression estimator
    import sys
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: UserWarning: Coordinate descent with no regularization may lead to unexpected results and is discouraged.
    positive)
```



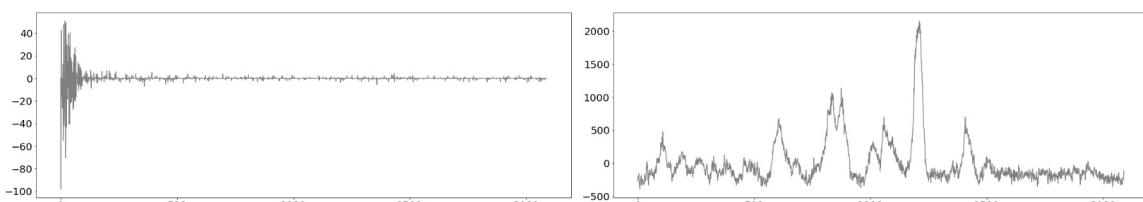
===== k = 0.3 , L1 = 0.001 =====



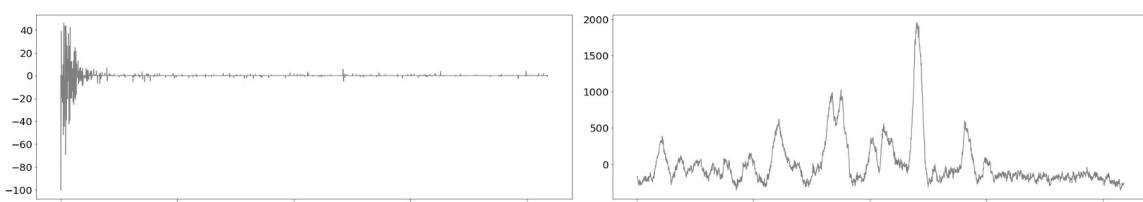
===== k = 0.3 , L1 = 0.01 =====



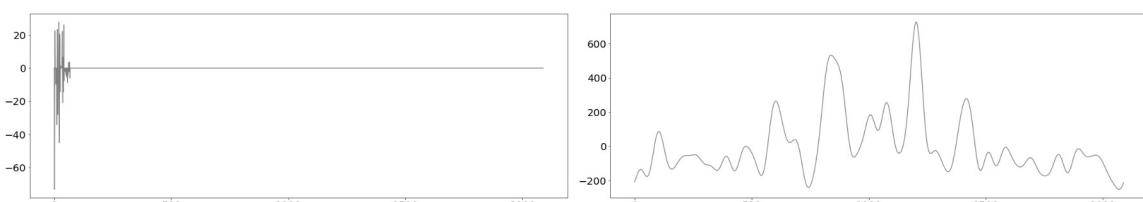
===== $k = 0.3$, $L1 = 0.1$ =====



===== $k = 0.3$, $L1 = 1$ =====



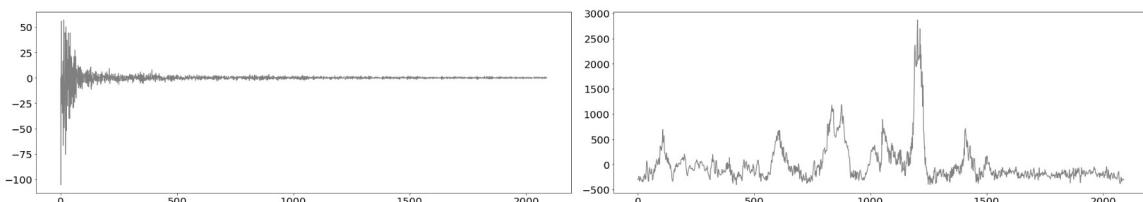
===== $k = 0.3$, $L1 = 10$ =====



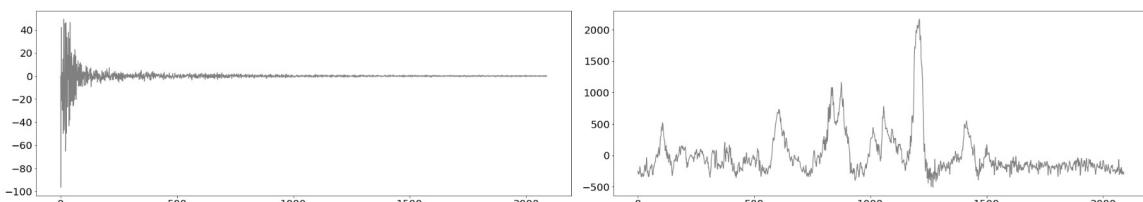
===== $k = 0.4$, $L1 = 0$ =====

```
C:\Users\eleve\Anaconda3\lib\site-packages\ipykernel_launcher.py:7: UserWarning: With alpha=0, this algorithm does not converge well. You are advised to use the LinearRegression estimator
```

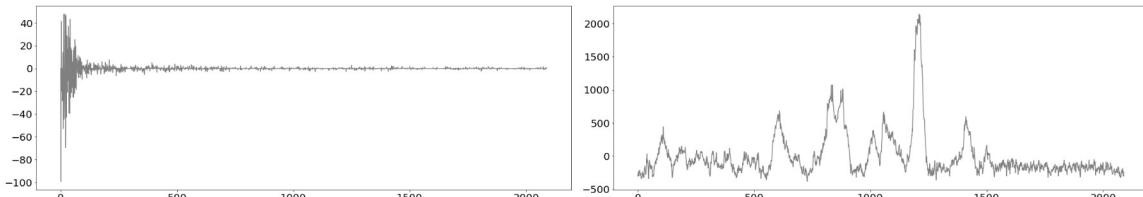
```
    import sys
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: UserWarning: Coordinate descent with no regularization may lead to unexpected results and is discouraged.
    positive)
```



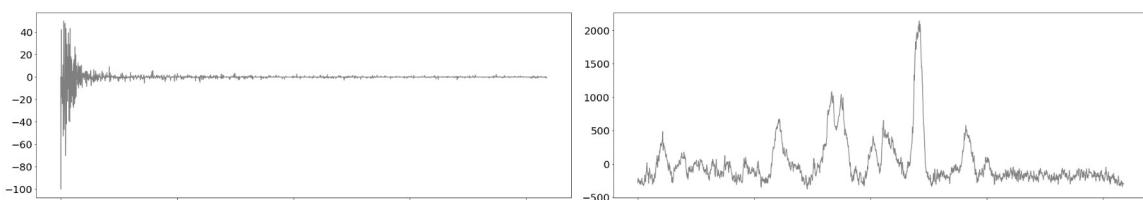
===== $k = 0.4$, $L1 = 0.001$ =====



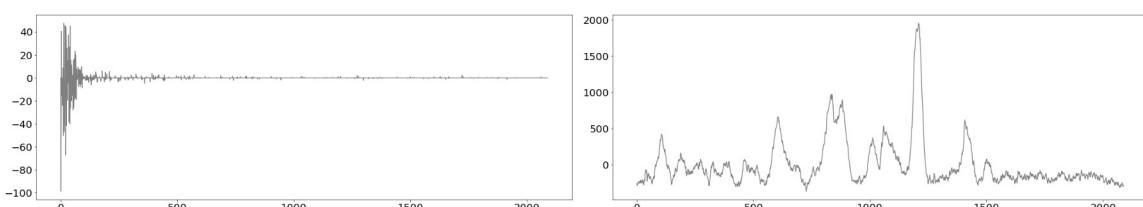
===== $k = 0.4$, $L1 = 0.01$ =====



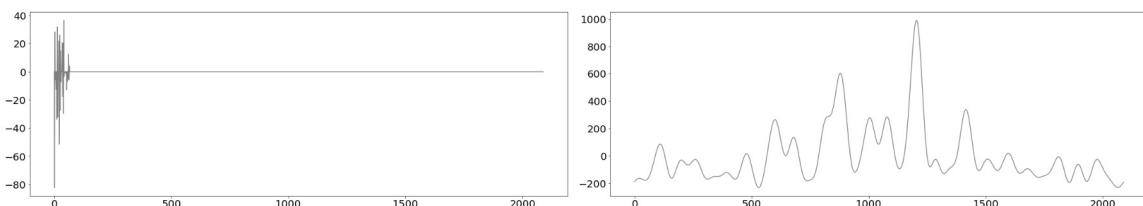
===== $k = 0.4$, $L_1 = 0.1$ =====



===== $k = 0.4$, $L_1 = 1$ =====



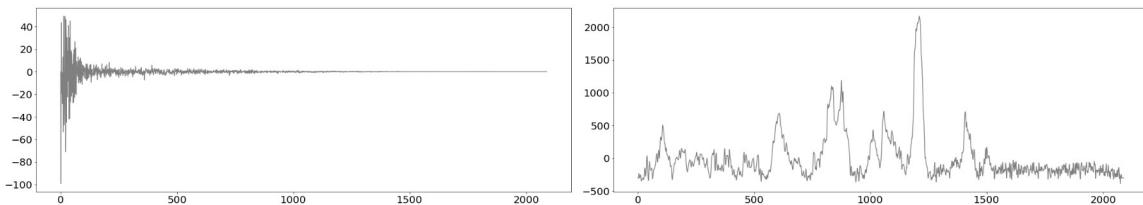
===== $k = 0.4$, $L_1 = 10$ =====



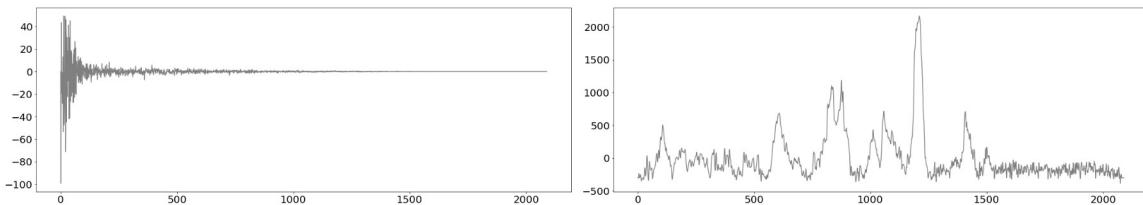
===== $k = 1.1$, $L_1 = 0$ =====

```
C:\Users\eleve\Anaconda3\lib\site-packages\ipykernel_launcher.py:7: UserWarning: With alpha=0, this algorithm does not converge well. You are advised to use the LinearRegression estimator
```

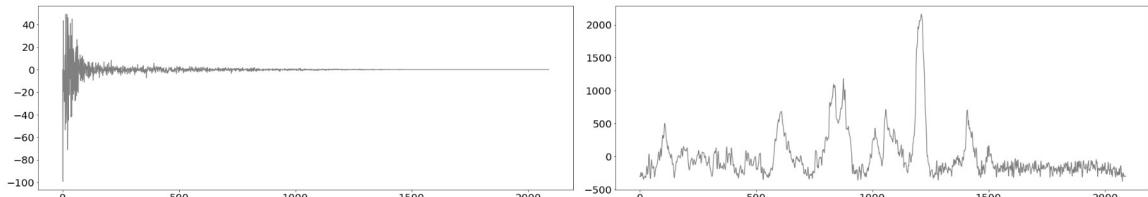
```
    import sys
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: UserWarning: Coordinate descent with no regularization may lead to unexpected results and is discouraged.
    positive)
```



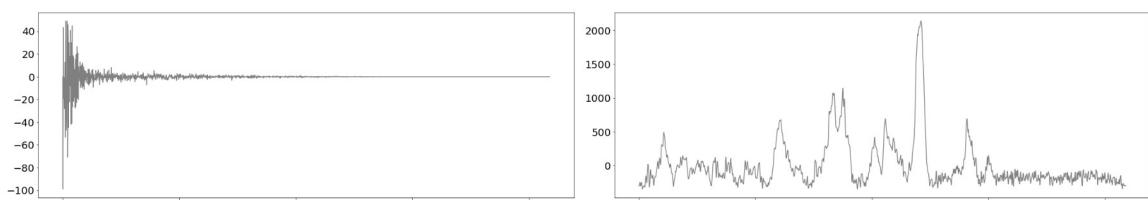
===== $k = 1.1$, $L_1 = 0.001$ =====



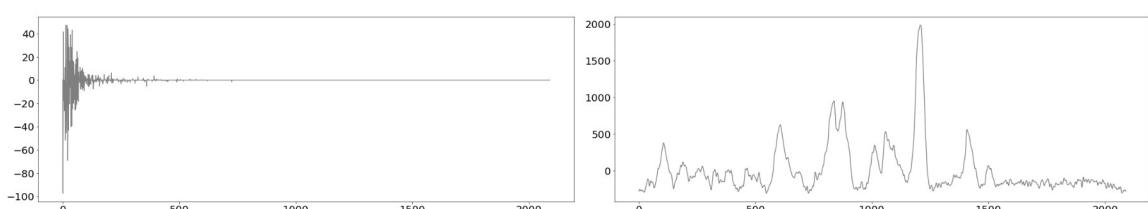
===== $k = 1.1$, $L_1 = 0.01$ =====



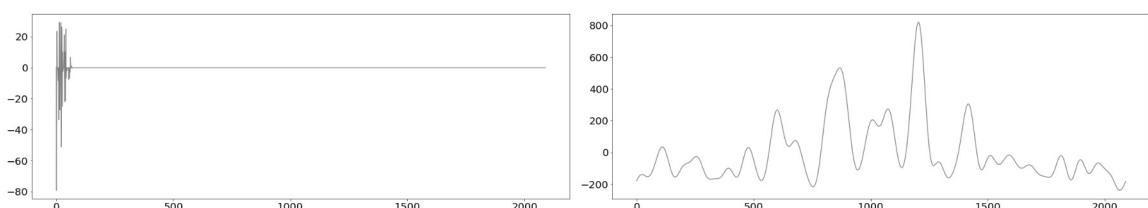
===== $k = 1.1$, $L_1 = 0.1$ =====



===== $k = 1.1$, $L_1 = 1$ =====



===== $k = 1.1$, $L_1 = 10$ =====



```
In [29]: import scipy

def all_samples_CS (X, y, k, L1, t='dft'):

    print('===== k =', k, ', L1 =', L1, ' =====')

    Z = np.zeros(X.shape)
    Xr = np.zeros(X.shape)

    if (k > 1.0): # when k > 1.0, return original signal directly
        Xr = X
        Z = X
    else:
        for i in range(X.shape[0]):
            x = X[i,:].ravel().tolist()[0] # get the first sample
            xr, z = one_sample_CS(x, k, L1, t)
            Z[i,:] = z
            Xr[i,:] = xr[:,0]

    fig, ax = plt.subplots(nrows=1, ncols=3, figsize=(20,4))

    from sklearn.decomposition import PCA

    pca = PCA(n_components=2) # keep the first 2 components
    Z_pca = pca.fit_transform(Z)
    # plotComponents2D(Z_pca, y, labels, use_markers = False, ax=ax[0])
    # ax[0].title.set_text('2D-Visualization')

    pca = PCA(n_components=2) # keep the first 2 components
    Xr_pca = pca.fit_transform(Xr)
    plotComponents2D(Xr_pca, y, labels, use_markers = False, ax=ax[0])
    ax[0].title.set_text('2D Visualization')

    Xc1s = []
    Xc2s = []
    title = ''
    for c in set(y):
        # print(Xr.shape, Z.shape, Xr_pca.shape, Z_pca.shape, y.shape)
        Xc = Xr_pca[y == c]
        yc = y[y == c]
        Xc1s.append(list(np.asarray(Xc[:,0]).reshape(1,-1)[0])) # First PC of Class c
        Xc2s.append(list(np.asarray(Xc[:,1]).reshape(1,-1)[0])) # Second PC of Class c

    #### ANOVA ####

    ax[1].title.set_text('PC1')
    ax[1].boxplot(Xc1s, notch=False) # plot 1st PC of all classes
    f,p= scipy.stats.f_oneway(Xc1s[0], Xc1s[1]) # equal to ttest_ind() in case of 2 groups
    print("f={},p={}".format(f,p))
    # dict_anova_p1[idx] = p

    ax[2].title.set_text('PC2')
    ax[2].boxplot(Xc2s, notch=False)
    f,p= scipy.stats.f_oneway(Xc2s[0], Xc2s[1])
    print("f={},p={}".format(f,p))

    #### MANOVA ####

    import statsmodels
    from statsmodels.base.model import Model
    from statsmodels.multivariate.manova import MANOVA

    df = pd.DataFrame({'PC1':Xr_pca[:,0], 'PC2':Xr_pca[:,1], 'y':y})
```

NOTE:

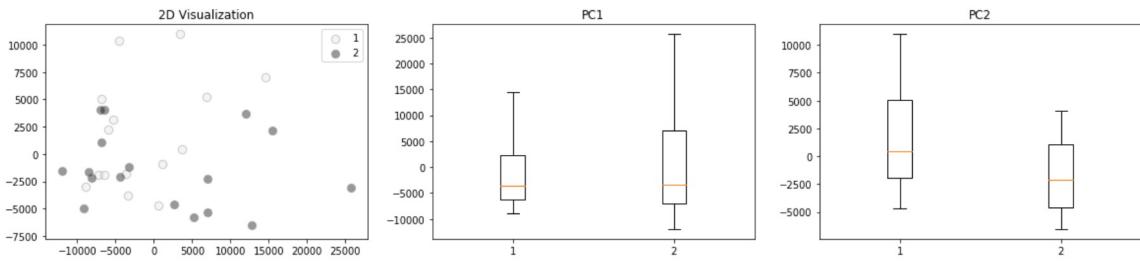
```
fig, axs = plt.subplots(3) returns a 1D array of subplots.  
fig, axs = plt.subplots(3, 2) returns a 2D array of subplots.
```

```
In [35]: result = {}

for k in [0.1, 0.2, 0.3, 0.4]:
    for L1 in [0.001, 0.01, 0.1, 1, 10]:
        acc, manova = all_samples_CS (X, y, k=k, L1=L1)
        result['k = ' + str(k) + 'L1 = ' + str(L1)] = (acc, manova)
```

```
===== k = 0.1 , L1 = 0.001 =====
f=0.7493537517132374,p=0.3935511106932027
f=4.77386574367189,p=0.036840699804596234
['PC1', 'PC2']
['Intercept', 'y']
    Multivariate linear model
=====
-----  
Intercept      Value  Num DF  Den DF F Value Pr > F  
-----  
Wilks' lambda 0.8516 2.0000 29.0000 2.5275 0.0973  
Pillai's trace 0.1484 2.0000 29.0000 2.5275 0.0973  
Hotelling-Lawley trace 0.1743 2.0000 29.0000 2.5275 0.0973  
Roy's greatest root 0.1743 2.0000 29.0000 2.5275 0.0973  
-----  
-----  
y      Value  Num DF  Den DF F Value Pr > F  
-----  
Wilks' lambda 0.8383 2.0000 29.0000 2.7959 0.0776  
Pillai's trace 0.1617 2.0000 29.0000 2.7959 0.0776  
Hotelling-Lawley trace 0.1928 2.0000 29.0000 2.7959 0.0776  
Roy's greatest root 0.1928 2.0000 29.0000 2.7959 0.0776  
=====
```

SVC ACC: 0.7827380952380952

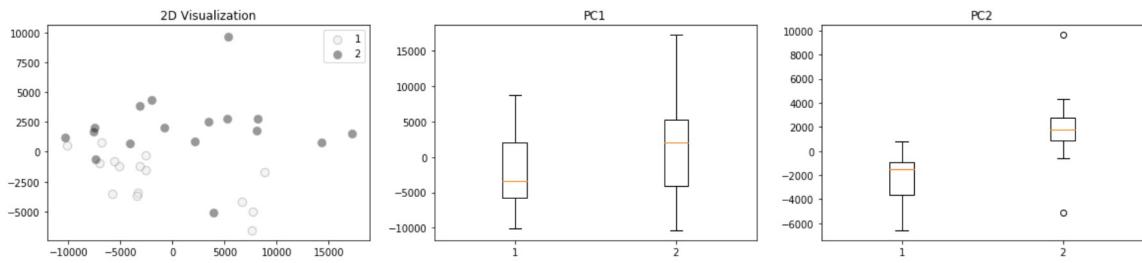


```
===== k = 0.1 , L1 = 0.01 =====
```

```
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 980.9499725346309, tolerance: 782.986769950874
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 995.5335269016138, tolerance: 512.2348822399686
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 1543.764280724276, tolerance: 1250.5539716828796
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 1671.1089656092902, tolerance: 1041.9075147361805
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 1192.4934233118297, tolerance: 892.2243608059954
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 1347.4624049441086, tolerance: 845.7480385807359
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 1172.0226256172343, tolerance: 813.8664765105298
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 2188.622850741333, tolerance: 1209.561852475047
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 1739.7490038627252, tolerance: 1047.1934162250868
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 962.6899210992066, tolerance: 569.8125551071149
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 1394.6425179301575, tolerance: 792.0628916975884
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 1804.9471506161372, tolerance: 1075.5320194624458
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 1998.4865673491176, tolerance: 1595.4903001871337
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 2632.7692895190144, tolerance: 2083.5362398239154
```

```
f=1.5432129009089282,p=0.22376380234100307
f=20.856841878700727,p=7.890254781126888e-05
['PC1', 'PC2']
['Intercept', 'y']
    Multivariate linear model
=====
-----
Intercept      Value  Num DF  Den DF F Value Pr > F
-----
Wilks' lambda 0.5659 2.0000 29.0000 11.1226 0.0003
Pillai's trace 0.4341 2.0000 29.0000 11.1226 0.0003
Hotelling-Lawley trace 0.7671 2.0000 29.0000 11.1226 0.0003
Roy's greatest root 0.7671 2.0000 29.0000 11.1226 0.0003
-----
Y      Value  Num DF  Den DF F Value Pr > F
-----
Wilks' lambda 0.5410 2.0000 29.0000 12.3038 0.0001
Pillai's trace 0.4590 2.0000 29.0000 12.3038 0.0001
Hotelling-Lawley trace 0.8485 2.0000 29.0000 12.3038 0.0001
Roy's greatest root 0.8485 2.0000 29.0000 12.3038 0.0001
=====
```

SVC ACC: 0.8809523809523809



===== k = 0.1 , L1 = 0.1 =====

```
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 685.1144116934224, tolerance: 504.20847236725206
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 676.217336259386, tolerance: 636.8756065497557
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 1234.824955222815, tolerance: 682.6001763826207
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 568.9776100222489, tolerance: 564.8443634706733
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 2487.825777561981, tolerance: 1168.2610343346375
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 664.2504767014262, tolerance: 433.00811377090014
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 1112.6569450264474, tolerance: 934.8281148181023
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 3758.735523306589, tolerance: 2424.9688982482103
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 2094.97137461006, tolerance: 1637.9793170312712
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 2562.312307470257, tolerance: 1926.7834575865627
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 3811.982269997713, tolerance: 1446.8529534519632
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 2242.861691395376, tolerance: 1480.2810270315763
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 2480.4866273775333, tolerance: 1220.2029818292021
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 3491.6832710325198, tolerance: 1761.9056237072994
```

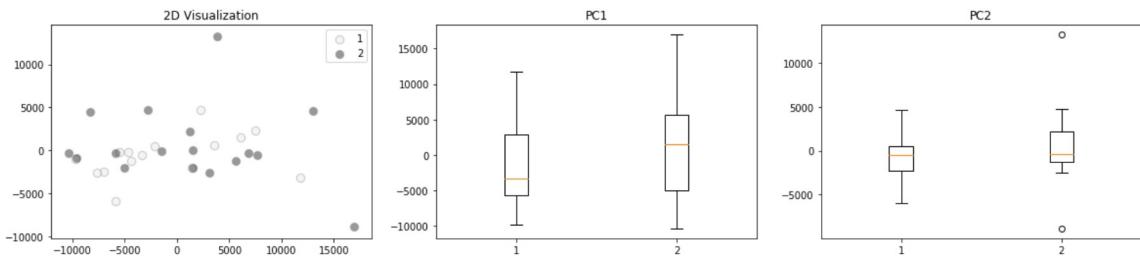
```
f=0.7941234480973275, p=0.3799478592173795
f=0.8744356668595947, p=0.3571968565216328
['PC1', 'PC2']
['Intercept', 'y']
    Multivariate linear model
=====
-----
```

Intercept	Value	Num DF	Den DF	F Value	Pr > F
Wilks' lambda	0.9508	2.0000	29.0000	0.7498	0.4814
Pillai's trace	0.0492	2.0000	29.0000	0.7498	0.4814
Hotelling-Lawley trace	0.0517	2.0000	29.0000	0.7498	0.4814
Roy's greatest root	0.0517	2.0000	29.0000	0.7498	0.4814

y	Value	Num DF	Den DF	F Value	Pr > F
Wilks' lambda	0.9459	2.0000	29.0000	0.8295	0.4464
Pillai's trace	0.0541	2.0000	29.0000	0.8295	0.4464
Hotelling-Lawley trace	0.0572	2.0000	29.0000	0.8295	0.4464
Roy's greatest root	0.0572	2.0000	29.0000	0.8295	0.4464

=====

SVC ACC: 0.7638888888888888



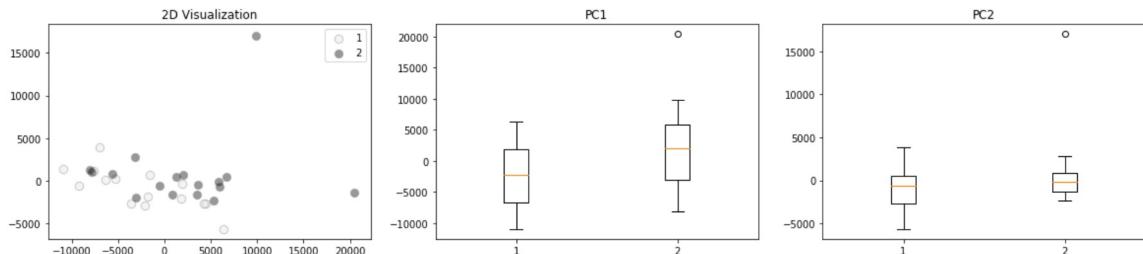
```
===== k = 0.1 , L1 = 1 =====
f=4.2966575063948715, p=0.04687407181404722
f=1.8430429403875888, p=0.18471983237259754
['PC1', 'PC2']
['Intercept', 'y']
    Multivariate linear model
=====
```

Intercept	Value	Num DF	Den DF	F Value	Pr > F
Wilks' lambda	0.8315	2.0000	29.0000	2.9391	0.0688
Pillai's trace	0.1685	2.0000	29.0000	2.9391	0.0688
Hotelling-Lawley trace	0.2027	2.0000	29.0000	2.9391	0.0688
Roy's greatest root	0.2027	2.0000	29.0000	2.9391	0.0688

y	Value	Num DF	Den DF	F Value	Pr > F
Wilks' lambda	0.8168	2.0000	29.0000	3.2513	0.0532
Pillai's trace	0.1832	2.0000	29.0000	3.2513	0.0532
Hotelling-Lawley trace	0.2242	2.0000	29.0000	3.2513	0.0532
Roy's greatest root	0.2242	2.0000	29.0000	3.2513	0.0532

=====

SVC ACC: 0.7941468253968255

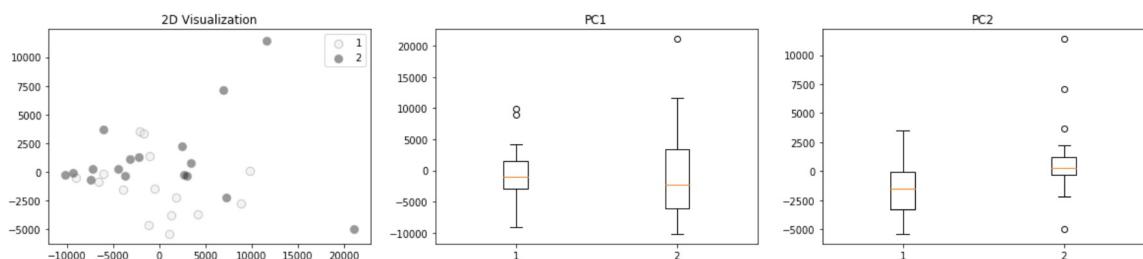


```
===== k = 0.1 , L1 = 10 =====
f=0.058248913655258854,p=0.8109273359030167
f=4.326647860749346,p=0.046161233069582044
['PC1', 'PC2']
['Intercept', 'y']
Multivariate linear model
=====
```

	Intercept	Value	Num DF	Den DF	F Value	Pr > F
Wilks' lambda	0.8829	2.0000	29.0000	1.9238	0.1642	
Pillai's trace	0.1171	2.0000	29.0000	1.9238	0.1642	
Hotelling-Lawley trace	0.1327	2.0000	29.0000	1.9238	0.1642	
Roy's greatest root	0.1327	2.0000	29.0000	1.9238	0.1642	

	Y	Value	Num DF	Den DF	F Value	Pr > F
Wilks' lambda	0.8720	2.0000	29.0000	2.1281	0.1373	
Pillai's trace	0.1280	2.0000	29.0000	2.1281	0.1373	
Hotelling-Lawley trace	0.1468	2.0000	29.0000	2.1281	0.1373	
Roy's greatest root	0.1468	2.0000	29.0000	2.1281	0.1373	

SVC ACC: 0.7271825396825398



```
===== k = 0.2 , L1 = 0.001 =====
f=2.658059154679605,p=0.11348254760962356
f=4.140535101357306,p=0.05078834064352125
['PC1', 'PC2']
['Intercept', 'y']
    Multivariate linear model
=====
-----
```

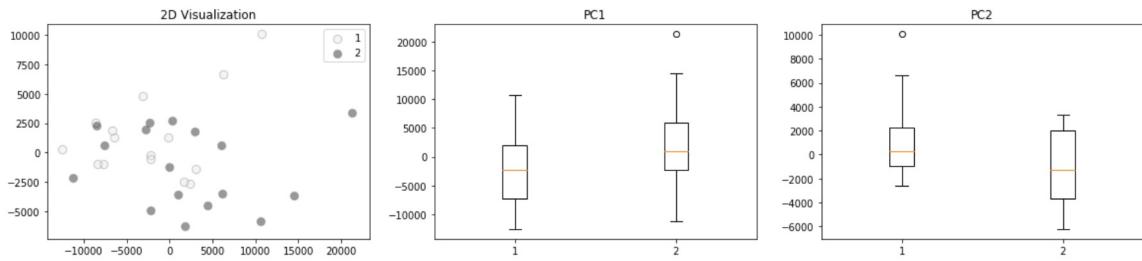
Intercept	Value	Num DF	Den DF	F Value	Pr > F
Wilks' lambda	0.8132	2.0000	29.0000	3.3318	0.0498
Pillai's trace	0.1868	2.0000	29.0000	3.3318	0.0498
Hotelling-Lawley trace	0.2298	2.0000	29.0000	3.3318	0.0498
Roy's greatest root	0.2298	2.0000	29.0000	3.3318	0.0498

```
-----
```

y	Value	Num DF	Den DF	F Value	Pr > F
Wilks' lambda	0.7973	2.0000	29.0000	3.6857	0.0375
Pillai's trace	0.2027	2.0000	29.0000	3.6857	0.0375
Hotelling-Lawley trace	0.2542	2.0000	29.0000	3.6857	0.0375
Roy's greatest root	0.2542	2.0000	29.0000	3.6857	0.0375

```
=====
```

SVC ACC: 0.9722222222222222



```
===== k = 0.2 , L1 = 0.01 =====
```

```
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 1648.3999843725676, tolerance: 1098.170854314675
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 2221.891225126621, tolerance: 1961.2839020810795
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 1659.7152402734075, tolerance: 1220.9177422381856
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 1453.6731511108164, tolerance: 1102.733587477971
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 5574.127337507545, tolerance: 4879.619859794887
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 3895.795615467818, tolerance: 3526.0940140691937
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 5609.0871545335085, tolerance: 5242.488328194646
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 6244.175570581737, tolerance: 5044.524647913703
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 12678.75547105898, tolerance: 10325.228789760262
    positive)
C:\Users\eleve\Anaconda3\lib\site-packages\sklearn\linear_model\coordinate_descent.py:475: ConvergenceWarning: Objective did not converge. You might want to increase the number of iterations. Duality gap: 5568.1748168475115, tolerance: 4482.609374454982
    positive)
```

```
f=1.724291902049198,p=0.19909735266364614
f=175.21170496720998,p=4.629902605206098e-14
['PC1', 'PC2']
['Intercept', 'y']
Multivariate linear model
=====

-----  

Intercept      Value  Num DF  Den DF F Value   Pr > F  

-----  

Wilks' lambda 0.1006 2.0000 29.0000 129.6202 0.0000
Pillai's trace 0.8994 2.0000 29.0000 129.6202 0.0000
Hotelling-Lawley trace 8.9393 2.0000 29.0000 129.6202 0.0000
Roy's greatest root 8.9393 2.0000 29.0000 129.6202 0.0000
-----  

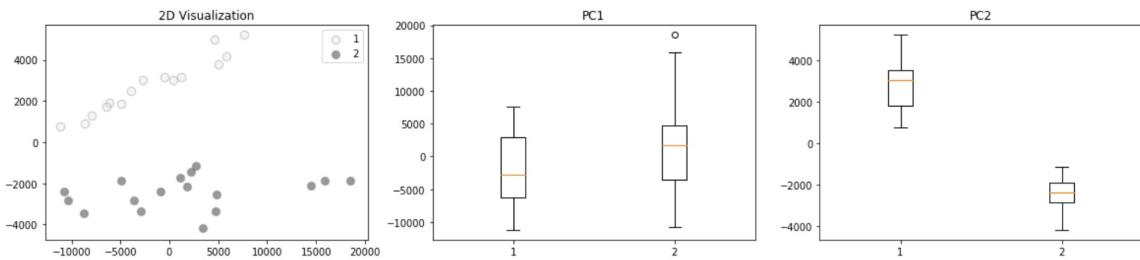
-----  

Y              Value  Num DF  Den DF F Value   Pr > F  

-----  

Wilks' lambda 0.0918 2.0000 29.0000 143.3866 0.0000
Pillai's trace 0.9082 2.0000 29.0000 143.3866 0.0000
Hotelling-Lawley trace 9.8887 2.0000 29.0000 143.3866 0.0000
Roy's greatest root 9.8887 2.0000 29.0000 143.3866 0.0000
=====
```

SVC ACC: 1.0



```
===== k = 0.2 , L1 = 0.1 =====
f=1.336008530334896,p=0.25686171105808797
f=97.89328168238266,p=5.85923666548548e-11
['PC1', 'PC2']
['Intercept', 'y']
Multivariate linear model
=====
```

```
-----  

Intercept      Value  Num DF  Den DF F Value   Pr > F  

-----  

Wilks' lambda 0.2081 2.0000 29.0000 55.1852 0.0000
Pillai's trace 0.7919 2.0000 29.0000 55.1852 0.0000
Hotelling-Lawley trace 3.8059 2.0000 29.0000 55.1852 0.0000
Roy's greatest root 3.8059 2.0000 29.0000 55.1852 0.0000
-----  

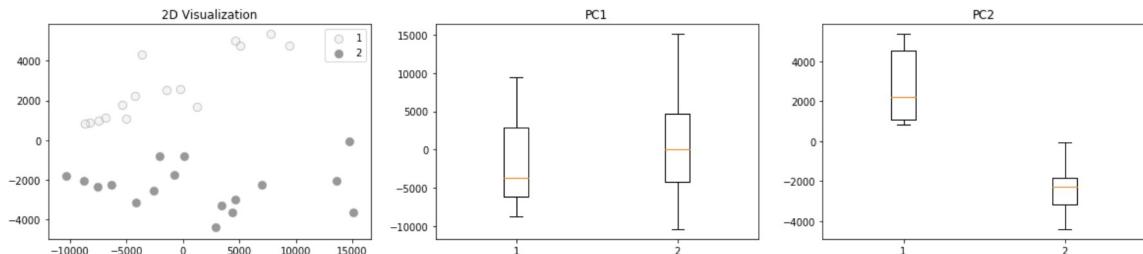
-----  

Y              Value  Num DF  Den DF F Value   Pr > F  

-----  

Wilks' lambda 0.1919 2.0000 29.0000 61.0462 0.0000
Pillai's trace 0.8081 2.0000 29.0000 61.0462 0.0000
Hotelling-Lawley trace 4.2101 2.0000 29.0000 61.0462 0.0000
Roy's greatest root 4.2101 2.0000 29.0000 61.0462 0.0000
=====
```

SVC ACC: 0.9444444444444444

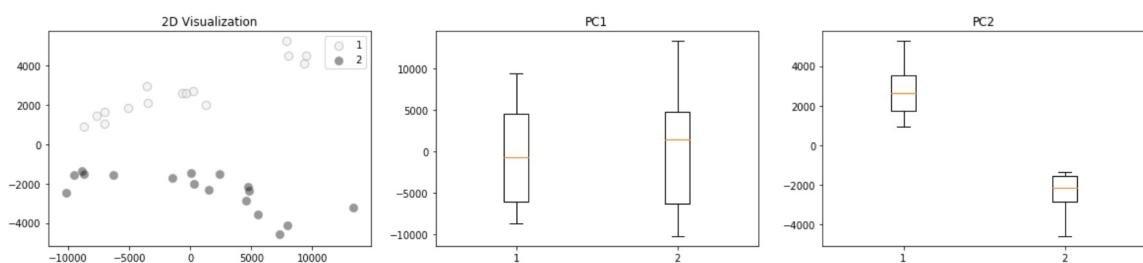


```
===== k = 0.2 , L1 = 1 =====
f=0.14773745500814037,p=0.7034181285033814
f=149.62170429158795,p=3.4530041297328493e-13
['PC1', 'PC2']
['Intercept', 'y']
Multivariate linear model
=====
```

```
-----
          Intercept      Value  Num DF  Den DF F Value Pr > F
-----
Wilks' lambda 0.1763 2.0000 29.0000 67.7464 0.0000
Pillai's trace 0.8237 2.0000 29.0000 67.7464 0.0000
Hotelling-Lawley trace 4.6722 2.0000 29.0000 67.7464 0.0000
Roy's greatest root 4.6722 2.0000 29.0000 67.7464 0.0000
-----
```

```
-----
          Y      Value  Num DF  Den DF F Value Pr > F
-----
Wilks' lambda 0.1621 2.0000 29.0000 74.9414 0.0000
Pillai's trace 0.8379 2.0000 29.0000 74.9414 0.0000
Hotelling-Lawley trace 5.1684 2.0000 29.0000 74.9414 0.0000
Roy's greatest root 5.1684 2.0000 29.0000 74.9414 0.0000
-----
```

SVC ACC: 0.9166666666666666



```
===== k = 0.2 , L1 = 10 =====
f=1.199788484435583,p=0.28208002935523246
f=63.75091540887328,p=6.523947170906379e-09
['PC1', 'PC2']
['Intercept', 'y']
Multivariate linear model
=====

-----  

Intercept      Value  Num DF  Den DF F Value Pr > F  

-----  

Wilks' lambda 0.3024 2.0000 29.0000 33.4496 0.0000
Pillai's trace 0.6976 2.0000 29.0000 33.4496 0.0000
Hotelling-Lawley trace 2.3069 2.0000 29.0000 33.4496 0.0000
Roy's greatest root 2.3069 2.0000 29.0000 33.4496 0.0000
-----  

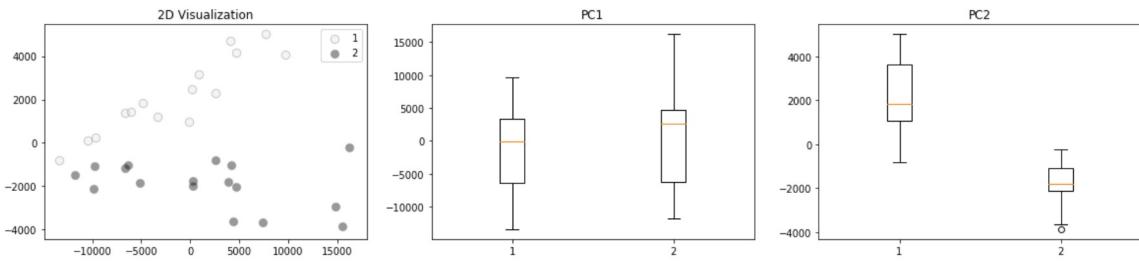
-----  

Y      Value  Num DF  Den DF F Value Pr > F  

-----  

Wilks' lambda 0.2815 2.0000 29.0000 37.0021 0.0000
Pillai's trace 0.7185 2.0000 29.0000 37.0021 0.0000
Hotelling-Lawley trace 2.5519 2.0000 29.0000 37.0021 0.0000
Roy's greatest root 2.5519 2.0000 29.0000 37.0021 0.0000
=====
```

SVC ACC: 0.8854166666666666



```
===== k = 0.3 , L1 = 0.001 =====
f=1.285999471719434,p=0.2657673068121966
f=78.41433845648044,p=7.170831586806164e-10
['PC1', 'PC2']
['Intercept', 'y']
Multivariate linear model
=====
```

```
-----  

Intercept      Value  Num DF  Den DF F Value Pr > F  

-----  

Wilks' lambda 0.2543 2.0000 29.0000 42.5255 0.0000
Pillai's trace 0.7457 2.0000 29.0000 42.5255 0.0000
Hotelling-Lawley trace 2.9328 2.0000 29.0000 42.5255 0.0000
Roy's greatest root 2.9328 2.0000 29.0000 42.5255 0.0000
-----  

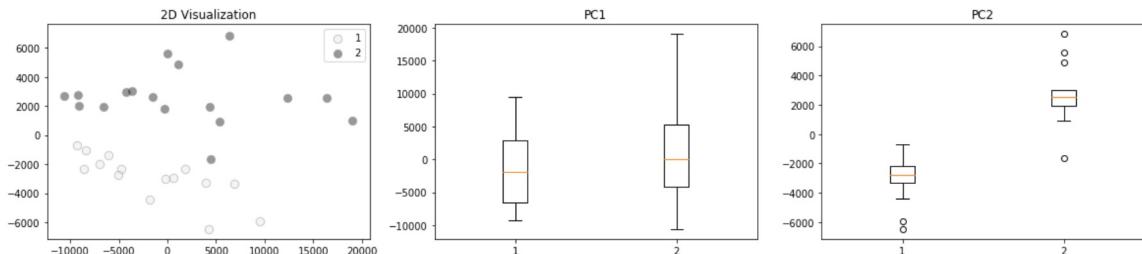
-----  

Y      Value  Num DF  Den DF F Value Pr > F  

-----  

Wilks' lambda 0.2356 2.0000 29.0000 47.0420 0.0000
Pillai's trace 0.7644 2.0000 29.0000 47.0420 0.0000
Hotelling-Lawley trace 3.2443 2.0000 29.0000 47.0420 0.0000
Roy's greatest root 3.2443 2.0000 29.0000 47.0420 0.0000
=====
```

SVC ACC: 1.0

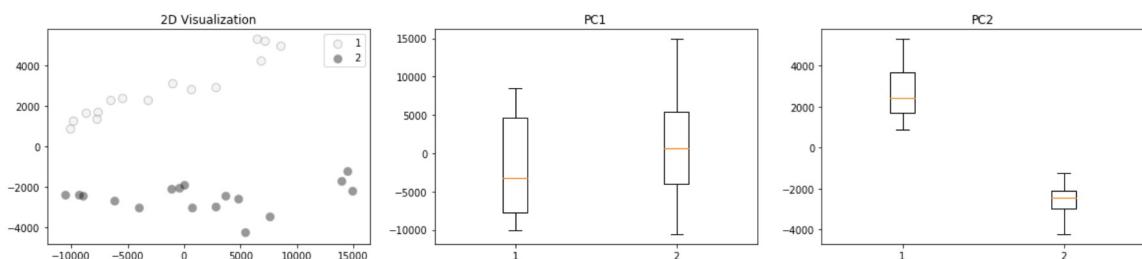


```
===== k = 0.3 , L1 = 0.01 =====
f=1.7212688966600238,p=0.1994808310215053
f=179.98582061057704,p=3.273186457707603e-14
['PC1', 'PC2']
['Intercept', 'y']
Multivariate linear model
=====
```

```
-----
          Intercept      Value  Num DF  Den DF F Value Pr > F
-----
Wilks' lambda 0.0971 2.0000 29.0000 134.8290 0.0000
Pillai's trace 0.9029 2.0000 29.0000 134.8290 0.0000
Hotelling-Lawley trace 9.2985 2.0000 29.0000 134.8290 0.0000
Roy's greatest root 9.2985 2.0000 29.0000 134.8290 0.0000
-----
```

```
-----
          Y      Value  Num DF  Den DF F Value Pr > F
-----
Wilks' lambda 0.0886 2.0000 29.0000 149.1486 0.0000
Pillai's trace 0.9114 2.0000 29.0000 149.1486 0.0000
Hotelling-Lawley trace 10.2861 2.0000 29.0000 149.1486 0.0000
Roy's greatest root 10.2861 2.0000 29.0000 149.1486 0.0000
-----
```

SVC ACC: 1.0

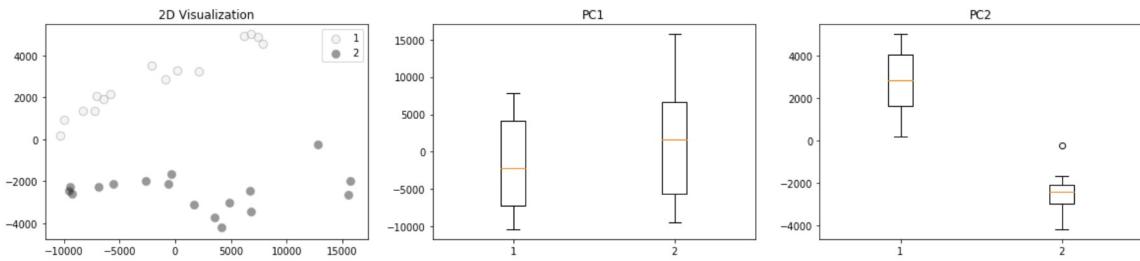


```
===== k = 0.3 , L1 = 0.1 =====
f=1.6747152781910628,p=0.20550221035108424
f=143.45245234577348,p=5.852117863984197e-13
['PC1', 'PC2']
['Intercept', 'y']
Multivariate linear model
=====
```

Intercept	Value	Num DF	Den DF	F Value	Pr > F
Wilks' lambda	0.1312	2.0000	29.0000	96.0464	0.0000
Pillai's trace	0.8688	2.0000	29.0000	96.0464	0.0000
Hotelling-Lawley trace	6.6239	2.0000	29.0000	96.0464	0.0000
Roy's greatest root	6.6239	2.0000	29.0000	96.0464	0.0000

y	Value	Num DF	Den DF	F Value	Pr > F
Wilks' lambda	0.1201	2.0000	29.0000	106.2470	0.0000
Pillai's trace	0.8799	2.0000	29.0000	106.2470	0.0000
Hotelling-Lawley trace	7.3274	2.0000	29.0000	106.2470	0.0000
Roy's greatest root	7.3274	2.0000	29.0000	106.2470	0.0000

SVC ACC: 1.0

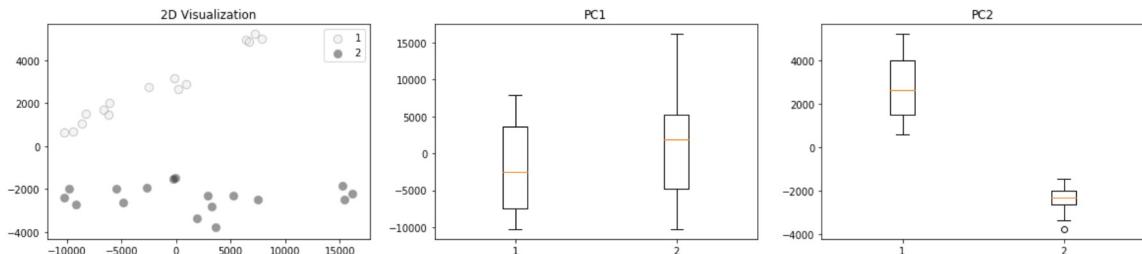


```
===== k = 0.3 , L1 = 1 =====
f=1.7804638023672994,p=0.19213287259429498
f=143.099357247646,p=6.0349511069829e-13
['PC1', 'PC2']
['Intercept', 'y']
Multivariate linear model
=====
```

Intercept	Value	Num DF	Den DF	F Value	Pr > F
Wilks' lambda	0.1281	2.0000	29.0000	98.6510	0.0000
Pillai's trace	0.8719	2.0000	29.0000	98.6510	0.0000
Hotelling-Lawley trace	6.8035	2.0000	29.0000	98.6510	0.0000
Roy's greatest root	6.8035	2.0000	29.0000	98.6510	0.0000

y	Value	Num DF	Den DF	F Value	Pr > F
Wilks' lambda	0.1173	2.0000	29.0000	109.1283	0.0000
Pillai's trace	0.8827	2.0000	29.0000	109.1283	0.0000
Hotelling-Lawley trace	7.5261	2.0000	29.0000	109.1283	0.0000
Roy's greatest root	7.5261	2.0000	29.0000	109.1283	0.0000

SVC ACC: 0.9444444444444444

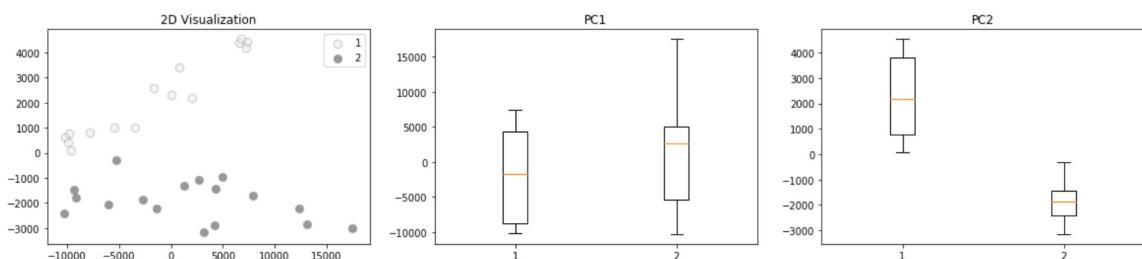


```
===== k = 0.3 , L1 = 10 =====
f=1.6129959507164962,p=0.21383388079884294
f=84.18839188733098,p=3.263762913987765e-10
['PC1', 'PC2']
['Intercept', 'y']
Multivariate linear model
=====
```

```
-----
          Intercept      Value  Num DF  Den DF F Value Pr > F
-----
Wilks' lambda 0.2290 2.0000 29.0000 48.8092 0.0000
Pillai's trace 0.7710 2.0000 29.0000 48.8092 0.0000
Hotelling-Lawley trace 3.3661 2.0000 29.0000 48.8092 0.0000
Roy's greatest root 3.3661 2.0000 29.0000 48.8092 0.0000
-----
```

```
-----
          Y      Value  Num DF  Den DF F Value Pr > F
-----
Wilks' lambda 0.2117 2.0000 29.0000 53.9930 0.0000
Pillai's trace 0.7883 2.0000 29.0000 53.9930 0.0000
Hotelling-Lawley trace 3.7237 2.0000 29.0000 53.9930 0.0000
Roy's greatest root 3.7237 2.0000 29.0000 53.9930 0.0000
-----
```

SVC ACC: 0.9444444444444444

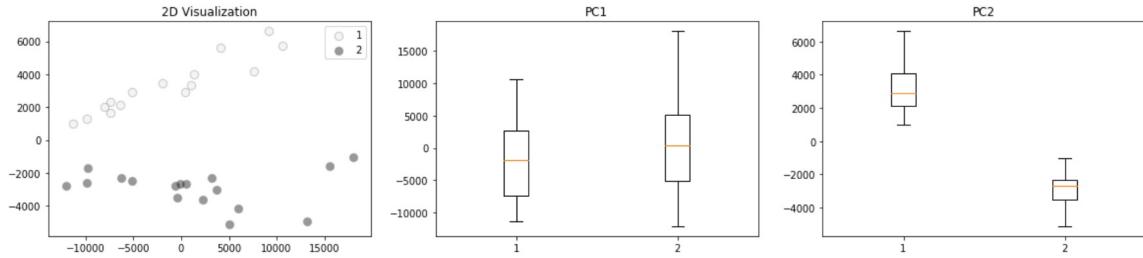


```
===== k = 0.4 , L1 = 0.001 =====
f=1.068507537376773,p=0.3095441681046772
f=154.54479015571607,p=2.2960341803484834e-13
['PC1', 'PC2']
['Intercept', 'y']
Multivariate linear model
=====
```

Intercept	Value	Num DF	Den DF	F Value	Pr > F
Wilks' lambda	0.1399	2.0000	29.0000	89.1614	0.0000
Pillai's trace	0.8601	2.0000	29.0000	89.1614	0.0000
Hotelling-Lawley trace	6.1491	2.0000	29.0000	89.1614	0.0000
Roy's greatest root	6.1491	2.0000	29.0000	89.1614	0.0000

y	Value	Num DF	Den DF	F Value	Pr > F
Wilks' lambda	0.1282	2.0000	29.0000	98.6308	0.0000
Pillai's trace	0.8718	2.0000	29.0000	98.6308	0.0000
Hotelling-Lawley trace	6.8021	2.0000	29.0000	98.6308	0.0000
Roy's greatest root	6.8021	2.0000	29.0000	98.6308	0.0000

SVC ACC: 0.9722222222222222

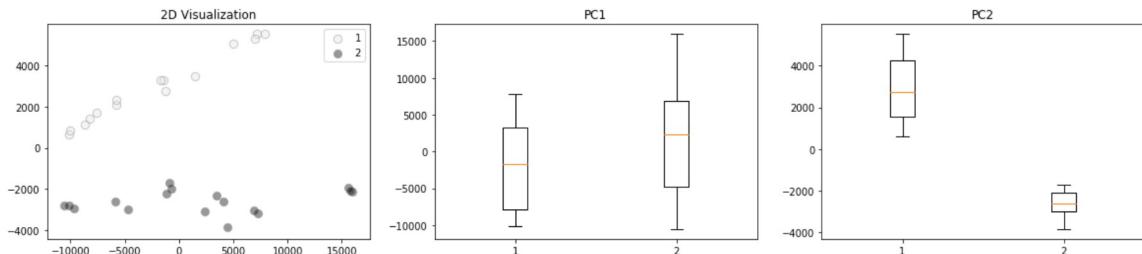


```
===== k = 0.4 , L1 = 0.01 =====
f=2.135114539818177,p=0.1543555513544617
f=157.54097638143037,p=1.8006900476903468e-13
['PC1', 'PC2']
['Intercept', 'y']
Multivariate linear model
=====
```

Intercept	Value	Num DF	Den DF	F Value	Pr > F
Wilks' lambda	0.1024	2.0000	29.0000	127.0484	0.0000
Pillai's trace	0.8976	2.0000	29.0000	127.0484	0.0000
Hotelling-Lawley trace	8.7620	2.0000	29.0000	127.0484	0.0000
Roy's greatest root	8.7620	2.0000	29.0000	127.0484	0.0000

y	Value	Num DF	Den DF	F Value	Pr > F
Wilks' lambda	0.0935	2.0000	29.0000	140.5417	0.0000
Pillai's trace	0.9065	2.0000	29.0000	140.5417	0.0000
Hotelling-Lawley trace	9.6925	2.0000	29.0000	140.5417	0.0000
Roy's greatest root	9.6925	2.0000	29.0000	140.5417	0.0000

SVC ACC: 1.0

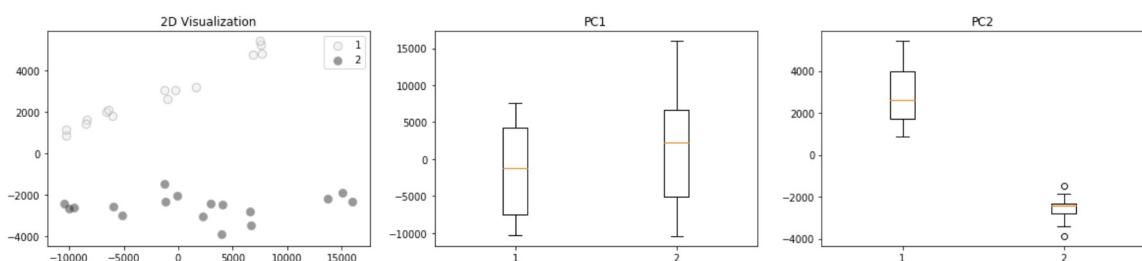


```
===== k = 0.4 , L1 = 0.1 =====
f=1.6376009151142177,p=0.21046333946071988
f=184.66081838039918,p=2.3485385856009768e-14
['PC1', 'PC2']
['Intercept', 'y']
Multivariate linear model
=====
```

```
-----
          Intercept      Value  Num DF  Den DF F Value Pr > F
-----
Wilks' lambda 0.0964 2.0000 29.0000 135.8551 0.0000
Pillai's trace 0.9036 2.0000 29.0000 135.8551 0.0000
Hotelling-Lawley trace 9.3693 2.0000 29.0000 135.8551 0.0000
Roy's greatest root 9.3693 2.0000 29.0000 135.8551 0.0000
-----
```

```
-----
          Y      Value  Num DF  Den DF F Value Pr > F
-----
Wilks' lambda 0.0880 2.0000 29.0000 150.2837 0.0000
Pillai's trace 0.9120 2.0000 29.0000 150.2837 0.0000
Hotelling-Lawley trace 10.3644 2.0000 29.0000 150.2837 0.0000
Roy's greatest root 10.3644 2.0000 29.0000 150.2837 0.0000
-----
```

SVC ACC: 1.0



```
===== k = 0.4 , L1 = 1 =====
f=1.747318565456052,p=0.196205730734803
f=156.90838627664965,p=1.8948707466927791e-13
['PC1', 'PC2']
['Intercept', 'y']
Multivariate linear model
=====
```

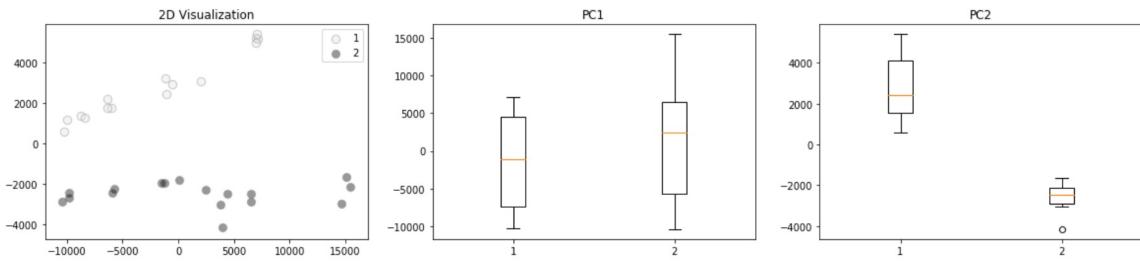
Intercept	Value	Num DF	Den DF	F Value	Pr > F
Wilks' lambda	0.1154	2.0000	29.0000	111.1749	0.0000
Pillai's trace	0.8846	2.0000	29.0000	111.1749	0.0000
Hotelling-Lawley trace	7.6672	2.0000	29.0000	111.1749	0.0000
Roy's greatest root	7.6672	2.0000	29.0000	111.1749	0.0000

=====

y	Value	Num DF	Den DF	F Value	Pr > F
Wilks' lambda	0.1055	2.0000	29.0000	122.9823	0.0000
Pillai's trace	0.8945	2.0000	29.0000	122.9823	0.0000
Hotelling-Lawley trace	8.4815	2.0000	29.0000	122.9823	0.0000
Roy's greatest root	8.4815	2.0000	29.0000	122.9823	0.0000

=====

SVC ACC: 1.0



```
===== k = 0.4 , L1 = 10 =====
f=0.6382368857090238,p=0.43062974045971825
f=115.33656099825497,p=8.46632373721209e-12
['PC1', 'PC2']
['Intercept', 'y']
Multivariate linear model
=====
```

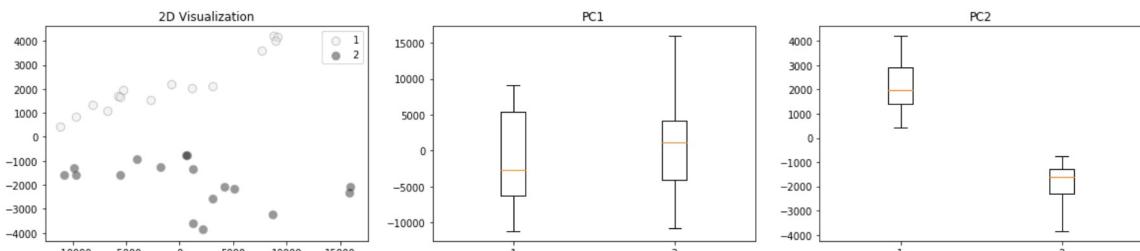
Intercept	Value	Num DF	Den DF	F Value	Pr > F
Wilks' lambda	0.2013	2.0000	29.0000	57.5217	0.0000
Pillai's trace	0.7987	2.0000	29.0000	57.5217	0.0000
Hotelling-Lawley trace	3.9670	2.0000	29.0000	57.5217	0.0000
Roy's greatest root	3.9670	2.0000	29.0000	57.5217	0.0000

=====

y	Value	Num DF	Den DF	F Value	Pr > F
Wilks' lambda	0.1856	2.0000	29.0000	63.6309	0.0000
Pillai's trace	0.8144	2.0000	29.0000	63.6309	0.0000
Hotelling-Lawley trace	4.3883	2.0000	29.0000	63.6309	0.0000
Roy's greatest root	4.3883	2.0000	29.0000	63.6309	0.0000

=====

SVC ACC: 1.0

**NOTE:**

No matter we use the sparse representation z in latent space, or the reconstructed data in the original space, the 2D visualizations after PCA are the same. Therefore, only the visualization of reconstructed x is shown.

When $k = 0.1$, the original discriminative structures in the data is lost (some samples from two classes begin to overlap), and the MANOVA under some λ test cannot reject the null-hypothesis H_0 .

$k = 0.2$ well preserves the discriminative structure. This result also supports the four-to-one rule.

```
In [34]: # Show K-fold ACC and MANOVA p-value in a table
```

```
from IPython.core.display import display, HTML

s = '<table>'

for key in result:
    s+= '<tr>'

    s+= '<td>' + key[:7] + '</td>'
    s+= '<td>' + key[7:] + '</td>'

    s+= '<td>' + str( round(result[key][0],3) ) + '</td>'
    s+= '<td>' + result[key][1].replace("Wilks' lambda", ", """)[:12] +
    '</td>'

    s+= '</tr>'

s += '</table>'
display(HTML(s))
```

k = 0.1	L1 = 0.001	0.783	0.0775615 Pi
k = 0.1	L1 = 0.01	0.881	0.000135201
k = 0.1	L1 = 0.1	0.764	0.446361 Pil
k = 0.1	L1 = 1	0.794	0.0532109 Pi
k = 0.1	L1 = 10	0.727	0.137285 Pil
k = 0.2	L1 = 0.001	0.972	0.0374761 Pi
k = 0.2	L1 = 0.01	1.0	9.2009e-16 P
k = 0.2	L1 = 0.1	0.944	4.03419e-11
k = 0.2	L1 = 1	0.917	3.48757e-12
k = 0.2	L1 = 10	0.885	1.04321e-08
k = 0.3	L1 = 0.001	1.0	7.88588e-10
k = 0.3	L1 = 0.01	1.0	5.47152e-16
k = 0.3	L1 = 0.1	1.0	4.49393e-14
k = 0.3	L1 = 1	0.944	3.19245e-14
k = 0.3	L1 = 10	0.944	1.67097e-10
k = 0.4	L1 = 0.001	0.972	1.15588e-13
k = 0.4	L1 = 0.01	1.0	1.19766e-15
k = 0.4	L1 = 0.1	1.0	4.9497e-16 P
k = 0.4	L1 = 1	1.0	6.84319e-15
k = 0.4	L1 = 10	1.0	2.47689e-11