Question 1

```
import numpy as np import pandas as pd import matplotlib.pyplot as plt from numpy import linalg as LA import time
```

```
In [2]:  # Load the data
    df1 = pd. read_csv('DailyReturn.csv')
    df1. describe()
```

ut[2]:		SPY	AAPL	MSFT	AMZN	TSLA	GOOGL	GOOG	FB	1
	count	60.000000	60.000000	60.000000	60.000000	60.000000	60.000000	60.000000	60.000000	60.0
	mean	0.000545	0.002621	0.000307	-0.000735	0.004213	-0.000149	-0.000193	-0.000239	0.0
	std	0.008858	0.016045	0.015962	0.016150	0.045304	0.015734	0.015441	0.020170	0.0
	min	-0.022303	-0.039264	-0.042323	-0.028955	-0.119903	-0.045876	-0.046830	-0.050515	-0.0
	25%	-0.003467	-0.005858	-0.007011	-0.012032	-0.028789	-0.006252	-0.006575	-0.011258	-0.0
	50%	0.000877	0.001231	0.000594	-0.001731	0.006190	-0.000209	0.000146	-0.001346	0.0
	75%	0.005470	0.015612	0.009943	0.007864	0.033007	0.008939	0.009319	0.015175	0.07
	max	0.020685	0.035446	0.042114	0.041437	0.135317	0.049595	0.048367	0.040123	0.17

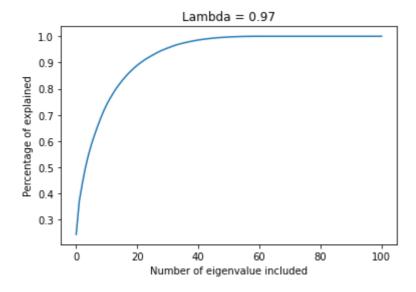
8 rows × 101 columns

```
# Write the function to get the exponentially weighted covariance matrix
def ewm (dataFrame, expect_weight, lamda):
    weight = []
    sum_weight = 0
    for i in range (1, 61):
        weight. append ((1-1 \text{amda})*1 \text{amda}**(i-1))
        sum_weight = sum_weight + weight[i-1]
    for i in range (0, 60):
        expect_weight.append(weight[i] / sum_weight)
    cov_matrix = np. zeros([101, 101])
    for i in range (0,60):
        for j in range (1, 102):
            dataFrame.iloc[i, j] = dataFrame.iloc[i, j] - np. mean(dataFrame.iloc[:, j])
    for i in range (0,101):
        for j in range (0, 101):
            #cov matrix.iloc[i,j] = sum( np.dot( np.dot(expect weight, dataFrame.iloc[
            temp = expect_weight * dataFrame.iloc[:,i+1]
            cov_matrix[i, j] = np. dot(temp, dataFrame.iloc[:, j+1])
```

return cov_matrix

```
In [4]:
          \# Lambda = 0.97
          expect_weight_1 = []
          cov_matrix_1 = ewm(df1, expect_weight_1, 0.97)
          value_1, vector_1 = LA. eigh(cov_matrix_1)
          value 1 = value 1[::-1]
          total 1 = 0
          for i in range(len(value_1)):
              total_1 = total_1 + value_1[i]
          cumu var 1 = []
          k = 1
          for i in range(len(value_1)):
              sum 1 = 0
              for j in range(k):
                  sum 1 = sum 1 + value 1[j]
              cumu_var_1. append(sum_1 / total_1)
              k = k + 1
          fig, ax = plt.subplots()
          ax. set_title("Lambda = 0.97")
          ax. set_xlabel('Number of eigenvalue included')
          ax. set_ylabel('Percentage of explained')
          ax. plot (cumu_var_1)
```

Out[4]: [<matplotlib.lines.Line2D at 0x17c07f8d1f0>]



```
In [5]: # Lambda = 0.7
    expect_weight_2 = []
    cov_matrix_2 = ewm(df1, expect_weight_2, 0.7)

value_2, vector_2 = LA. eigh(cov_matrix_2)
    value_2 = value_2[::-1]

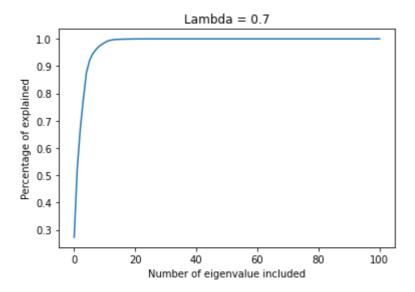
total_2 = 0
    for i in range(len(value_2)):
        total_2 = total_2 + value_2[i]

cumu_var_2 = []
    k = 1
```

```
for i in range(len(value_2)):
    sum_2 = 0
    for j in range(k):
        sum_2 = sum_2 + value_2[j]
    cumu_var_2. append(sum_2 / total_2)
    k = k + 1

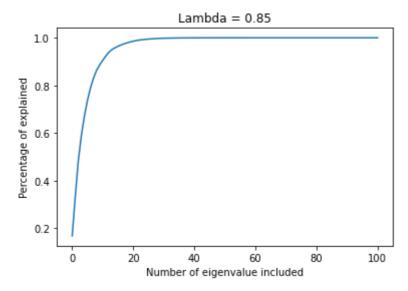
fig, ax = plt. subplots()
ax. set_title("Lambda = 0.7")
ax. set_xlabel('Number of eigenvalue included')
ax. set_ylabel('Percentage of explained')
ax. plot(cumu_var_2)
```

Out[5]: [<matplotlib.lines.Line2D at Ox17c0a8b6550>]



```
In [6]:
          # Lambda = 0.85
          expect_weight_3 = []
          cov_matrix_3 = ewm(df1, expect_weight_3, 0.85)
          value_3, vector_3 = LA. eigh(cov_matrix_3)
          value_3 = value_3[::-1]
          total 3 = 0
          for i in range(len(value 3)):
              total_3 = total_3 + value_3[i]
          cumu_var_3 = []
          k = 1
          for i in range(len(value_3)):
              sum_3 = 0
              for j in range(k):
                  sum 3 = sum 3 + value 3[j]
              cumu var 3. append(sum 3 / total 3)
              k = k + 1
          fig, ax = plt. subplots()
          ax. set_title("Lambda = 0.85")
          ax. set xlabel('Number of eigenvalue included')
          ax. set_ylabel('Percentage of explained')
          ax. plot (cumu_var_3)
```

Out[6]: [<matplotlib.lines.Line2D at 0x17c0a915ca0>]

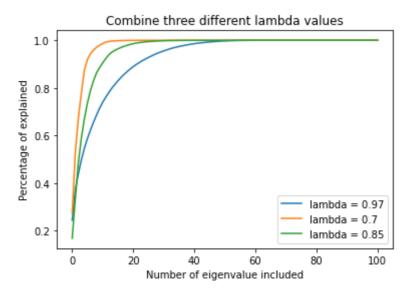


```
fig, ax = plt. subplots()
ax. set_title("Combine three different lambda values")
ax. set_xlabel('Number of eigenvalue included')
ax. set_ylabel('Percentage of explained')

ax. plot(cumu_var_1, label = 'lambda = 0.97')
ax. plot(cumu_var_2, label = 'lambda = 0.7')
ax. plot(cumu_var_3, label = 'lambda = 0.85')

ax. legend()
```

Out[7]: <matplotlib.legend.Legend at 0x17c0a8e68e0>



Question 2

```
In [8]:  # non-psd correlation matrix that is 500x500
    n = 500
    sigma = np. full((n,n), 0.9)
    for i in range(n):
        sigma[i,i] = 1.0
    sigma[1,2] = 0.7357
    sigma[2,1] = 0.7357
```

In [9]:

```
# determined whether the matix is psd
def is_psd(x):
   return np. all(np. linalg. eigvals(x) >= -1e-8)
```

```
In [10]: # Near PSD function
def near_psd(A):
    n = A.shape[0]
    eigval, eigvec = LA.eigh(A)
    val = np.matrix(np.maximum(eigval, 0))
    vec = np.matrix(eigvec)

    T = 1/(np.multiply(vec, vec) * val.T)
    T = np.matrix(np.sqrt(np.diag(np.array(T).reshape((n))))) #already take the squa
    B = T * vec * np.diag(np.array(np.sqrt(val)).reshape((n)))
    out = B * B.T

    return out
```

```
In [11]:
           # Higham PSD function
           def _getAplus(A):
               eigval, eigvec = LA. eigh(A)
               Q = np. matrix(eigvec)
               xdiag = np. matrix(np. diag(np. maximum(eigval, 0)))
               return eigvec @ xdiag @ eigvec.T
           def getPs(A, W=None):
               W05 = np. matrix(W**.5)
               iW = W05. I
               return iW @ _getAplus(W05 @ A @ W05) @ iW
           def _getPu(A, W=None):
               Aret = A. copy()
               for i in range(0, A. shape[0]):
                   Aret[i, i] = 1
               return Aret
           def _wgtNorm(A, W = None):
               W05 = np. sqrt(W)
               W05 = W05 @ A @ W05
               return (W05 * W05). sum()
           def higham nearestPSD(pc, W = None, epsilon = 1e-9, maxIter = 100, tol = 1e-9):
               n = pc. shape[0]
               if W == None:
                   W = np. identity(n)
               deltaS = np. zeros((n, n))
               Yk = pc. copy()
               norm1 = 99999999
               i = 1
               while i <= maxIter:
                   Rk = Yk - deltaS
                   Xk = _{getPs(Rk, W)}
                   deltaS = Xk - Rk
                   Yk = getPu(Xk, W)
                   norm = \_wgtNorm(Yk - pc, W)
```

w, v = LA. eigh(Yk)

```
minEigVal = np. min(w)

if ((norm - norml) < tol) and (minEigVal > -epsilon):
    break

norml = norm
    i = i + 1

if i < maxIter:
    print("Converged in %d iterations. \n" % i)
else:
    print("Converged failed after %d iterations. \n" % (i-1))

return Yk</pre>
```

```
# Cholesky PSD Function
def chol_psd(root, a):
    n = a. shape[0]
    for j in range(n):
        s = 0.0
        if j > 0:
            s = root[j, :j] @ root[j, :j].T
        temp = a[j, j] - s
        if -1e-8 \le temp \le 0:
            temp = 0.0
        root[j, j] = np. sqrt(temp);
        if root[j, j] == 0.0:
            continue
        ir = 1.0/root[j, j]
        for i in range (j+1, n):
                s = root[i, :j] @ root[j, :j].T
                root[i, j] = (a[i, j] - s) * ir
    return root
```

```
# Test whether the covariance is PSD after using near_psd function
out_near_psd = near_psd(sigma)
test_near_psd = is_psd(out_near_psd)
if test_near_psd == True:
    print("The covariance martix is PSD by using near PSD method.")
```

The covariance martix is PSD by using near PSD method.

```
# Test whether the covariance is PSD after using higham_nearestPSD function out_higham_nearestPSD = higham_nearestPSD(sigma, W = None, epsilon = 1e-9, maxIter = test_higham_nearestPSD = is_psd(out_higham_nearestPSD) if test_higham_nearestPSD == True:

print("The covariance martix is PSD by using Higham method.")
```

Converged in 26 iterations.

The covariance martix is PSD by using Higham method.

```
In [15]: # Frobenius norm function
```

```
def F_norm (cov, cov_psd):
    temp = cov - cov_psd
    return LA. norm(temp, 'fro')
```

```
# Calculate the Frobenius norm for both methods.

near_psd_Fnorm = F_norm(sigma, out_near_psd)
higham_psd_Fnorm = F_norm(sigma, out_higham_nearestPSD)

print("Frobenius Norm of near PSD method is %f" % near_psd_Fnorm)
print("Frobenius Norm of Higham PSD method is %f" %higham_psd_Fnorm)
```

Frobenius Norm of near PSD method is 0.627523 Frobenius Norm of Higham PSD method is 0.089648

```
In [17]:  # Calculate time consuming
    start1 = time.time()
    near_psd(sigma)
    end1 = time.time()
    near_psd_time = end1 - start1

    start2 = time.time()
    higham_nearestPSD(sigma, W = None, epsilon = 1e-9, maxIter = 100, tol = 1e-9)
    end2 = time.time()
    higham_psd_time = end2 - start2

    print("Near PSD method time consuming is %f" % near_psd_time)
    print("Near Higham method time consuming is %f" % higham_psd_time)
```

Converged in 26 iterations.

Near PSD method time consuming is 0.028581 Near Higham method time consuming is 0.938344

```
        Out[18]:
        Name
        Norm
        RunTime

        0
        nearPSD
        0.627523
        0.028581

        1
        Higham2002
        0.089648
        0.938344
```

Question 3

```
In [19]: # Cholesky PSD Function
    def chol_psd(root, a):
        n = a. shape[0]
    for j in range(n):
```

```
s = 0.0
        if j > 0:
            s = root[j, :j] @ root[j, :j].T
        temp = a[j, j] - s
        if -1e-8 \le temp \le 0:
            temp = 0.0
        root[j, j] = np. sqrt(temp);
        if root[j, j] == 0.0:
            continue
        ir = 1.0/root[j, j]
        for i in range (j+1, n):
                s = root[i, :j] @ root[j, :j].T
                root[i, j] = (a[i, j] - s) * ir
    return root
# Frobenius norm function
def F norm (cov, cov psd):
    temp = cov - cov_psd
    return LA.norm(temp, 'fro')
def getCor(cov):
    cov_diag = np. diag(cov)
    invSD = np. diag(np. divide(1, np. sqrt(cov_diag)))
    cor = invSD * cov * invSD
    return cor
# Standard Pearson correlation/variance
sp_corr = np. matrix(df1. corr())
sp_var = df1. var()
# Expoentially weighted lambda = 0.97
ew var = np. diag(cov matrix 1)
cov_matrix_1 = np. matrix(cov_matrix_1)
ew_corr = getCor(cov_matrix_1)
```

Combine these variance and correlation to form 4 different covariance matrix

```
In [24]:  # Combine Standard Peason's Variables and Standard Peason's Correlation PEARSON = combineVar(sp_var, sp_corr)
```

Combine Standard Peason's Variables and Exponentially Weighted's Correlation

```
EWMA COR PEARSON STD = combineVar(sp var, ew corr)
           # Combine Exponentially Weighted's Variables and Exponentially Weighted's Correlation
           EWMA = combineVar(ew_var, ew_corr)
           # Combine Exponentially Weighted's Variables and Standard Peason's Correlation
           PEARSON_COR_EWMA_STD = combineVar(ew_var, sp_corr)
In [28]:
           # Direct Simulation Function
           def simulateNormal(cov, nsim):
               if (cov. shape [0] != len(cov)):
                   raise exception ("covariance matrix is not square")
               n = cov. shape[0]
               root = np. zeros (cov. shape)
               root = chol_psd(root, cov)
               np. random. seed (1998)
               z = np. random. normal(size=(n, nsim))
               ans = root @ z
               return ans
           # PCA Simulation Function
           def simulate pca(cov, nsim, target):
               val, vec = LA. eigh(cov)
               tot = sum(val)
               va1 = va1[::-1]
               vec = vec[::-1]
               cumm val explained = np. cumsum(val) / tot
               i = 0
               for i in range(len(val)):
                   if cumm val explained[i] < target:</pre>
                       i += 1
                   else:
                       break
               val = val[0:i+1]
               vec = vec[:, :i+1]
               B = vec @ np. diag(np. sqrt(val))
               np. random. seed (1998)
               z = np. random. normal(size=(1en(val), nsim))
               return B @ z
           # Print Function
           matrixType = ["PEARSON", "EWMA COR PEARSON STD", "EWMA", "PEARSON COR EWMA STD"]
           simType = ["Full", "PCA=1", "PCA=0.75", "PCA=0.5"]
           matrix = []
           simulation = []
           runtimes = []
           norms = []
           i = 0
           nsim = 25000
```

```
for sim in simType:
    for mat in matrixType:
        global i
        matrix.append(mat)
        simulation. append(sim)
        elapse = 0
        if mat == "PEARSON":
            c = PEARSON
        elif mat == "EWMA_COR_PEARSON_STD":
            c = EWMA\_COR\_PEARSON\_STD
        elif mat == "EWMA":
            c = EWMA
        elif mat == "PEARSON COR EWMA STD":
            c = PEARSON\_COR\_EWMA\_STD
        if sim == 'Full':
            start = time. time()
            s = simulateNormal(c, nsim)
            end = time. time()
            elapse = end - start
        elif sim == 'PCA=1':
            start = time. time()
            s = simulate_pca(c, nsim, 1)
            end = time. time()
            elapse = end - start
        elif sim == 'PCA=0.75':
            start = time. time()
            s = simulate_pca(c, nsim, 0.75)
            end = time. time()
            elapse = end - start
        elif sim == 'PCA=0.5':
            start = time. time()
            s = simulate_pca(c, nsim, 0.5)
            end = time. time()
            elapse = end - start
        covar = np. cov(s)
        runtimes. append (elapse)
        norms.append(F norm(covar, c))
        i = i + 1
outTable = pd. DataFrame(list(zip(matrix, simulation, norms, runtimes,)), columns = ['
outTable
```

Out[30]:		Name	Simulation	Norm	RunTime
	0	PEARSON	Full	0.000188	0.065420
	1	EWMA_COR_PEARSON_STD	Full	0.000194	0.086290
	2	EWMA	Full	0.000185	0.086526
	3	PEARSON_COR_EWMA_STD	Full	0.000180	0.084338
	4	PEARSON	PCA=1	0.016373	0.044749
	5	EWMA_COR_PEARSON_STD	PCA=1	0.014891	0.044611
	6	EWMA	PCA=1	0.014138	0.046169
	7	PEARSON_COR_EWMA_STD	PCA=1	0.015429	0.040044
	8	PEARSON	PCA=0.75	0.016434	0.010963

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	Name	Simulation	Norm	RunTime
9	EWMA_COR_PEARSON_STD	PCA=0.75	0.014960	0.010965
10	EWMA	PCA=0.75	0.014188	0.012957
11	PEARSON_COR_EWMA_STD	PCA=0.75	0.015492	0.012956
12	PEARSON	PCA=0.5	0.016318	0.007499
13	EWMA_COR_PEARSON_STD	PCA=0.5	0.014772	0.007973
14	EWMA	PCA=0.5	0.014050	0.008970
15	PEARSON COR EWMA STD	PCA=0.5	0.015358	0.007549