R Outputs related to PCA for n < p

Dimension of the Data

```
> View(X)
> dim(X)
[1] 64 6830
```

Dimension of the Variance-Covariance Matrix

```
> c = cov(X)
> dim(c)
[1] 6830 6830
```

Eigenvalues of Var-Cov Matrix

```
= eigen(c)
>
> round(e$values,4)
   [1] 633.2156 352.9278 279.9189 183.0830 163.5573 149.0968
       122.2882 119.7912 112.1777
                                     91.7108
                                               89.2895
                                                         85.1300
                           71.0961
  [13]
        77.8269
                  75.1452
                                     68.4686
                                               67.0811
                                                         61.8224
  [19]
        61.5536
                  60.0242
                            58.3950
                                     54.9767
                                               53.5923
                                                         50.3407
  [25]
        50.1501
                  47.1681
                            45.1392
                                     43.8107
                                               43.2037
                                                         42.2685
  [31]
        40.7568
                  39.9030
                            37.2435
                                     36.8453
                                               35.5733
                                                         35.2031
  [37]
        34.4904
                  33.9733
                            32.2537
                                     31.7271
                                               31.0329
                                                         30.3893
  [43]
        29.7626
                  28.9381
                            28.3175
                                     26.7731
                                               26.4679
                                                         25.1793
  [49]
        23.2744
                  22.8368
                            22.0119
                                     20.8277
                                               20.1636
                                                         19.4607
  [55]
                            16.6965
                                     15.3630
                                               14.3497
        18.2677
                  17.7540
                                                         12.4190
  [61]
        10.4253
                   9.9400
                             8.9138
                                       0.0000
                                                0.0000
                                                          0.0000
                             0.0000
                                       0.0000
                                                0.0000
  [67]
         0.0000
                   0.0000
                                                          0.0000
  [73]
         0.0000
                   0.0000
                             0.0000
                                       0.0000
                                                0.0000
                                                          0.0000
  [79]
         0.0000
                   0.0000
                             0.0000
                                       0.0000
                                                0.0000
                                                          0.0000
  [85]
         0.0000
                   0.0000
                             0.0000
                                       0.0000
                                                0.0000
                                                          0.0000
  [91]
         0.0000
                   0.0000
                             0.0000
                                       0.0000
                                                0.0000
                                                          0.0000
  [97]
         0.0000
                   0.0000
                             0.0000
                                       0.0000
                                                0.0000
                                                          0.0000
 [103]
         0.0000
                   0.0000
                             0.0000
                                       0.0000
                                                0.0000
                                                          0.0000
 [109]
         0.0000
                   0.0000
                             0.0000
                                       0.0000
                                                0.0000
                                                          0.0000
 [115]
         0.0000
                   0.0000
                             0.0000
                                       0.0000
                                                0.0000
                                                          0.0000
 [121]
         0.0000
                   0.0000
                             0.0000
                                       0.0000
                                                0.0000
                                                          0.0000
 [127]
         0.0000
                   0.0000
                             0.0000
                                       0.0000
                                                0.0000
                                                          0.0000
 [133]
         0.0000
                   0.0000
                             0.0000
                                       0.0000
                                                0.0000
                                                          0.0000
 [139]
         0.0000
                   0.0000
                             0.0000
                                       0.0000
                                                0.0000
                                                          0.0000
 [145]
         0.0000
                   0.0000
                             0.0000
                                       0.0000
                                                0.0000
                                                          0.0000
```

Note that only first 63 eigenvalues are non-zero. The remaining (6830-63) eigenvalues are all equal to zero.

Rank of Var-Cov Matrix

```
> require(Matrix)
> r = rankMatrix(c)
> r[1]
[1] 63
```

Note that the rank of the variance — covariance matrix is 63, which is equal to the number of non-zero eigenvalues.