Factor Analysis on K-Medoid clusters Factor Analysis on 5th cluster In [1]: import pandas as pd from factor_analyzer import FactorAnalyzer from factor analyzer.factor analyzer import calculate_bartlett_sphericity, calculate_kmo import seaborn as sns import matplotlib.pyplot as plt #First the unnecessary columns are removed: df = pd.read excel('Asansol MRBQ Data.xlsx') df = df.drop(columns = ['TOT2']) df_2 = pd.read_excel('clustered data KMedoids.xlsx') clusters = list(df_2['Cluster']) df['Cluster'] = clusters df Out[1]: Q1 Q2 Q3 Q4 Q5 Q6 Q7 Q8 Q9 Q10 ... Q28 Q29 Q30 Q31 Q32 Q33 Q34 Q35 Q36 Cluster 4 3 476 2 ... 481 rows × 37 columns In [2]: # Only the rows belionging to the 5th cluster are filtered out. cluster_5 = df.loc[df['Cluster']==4] cluster 5 Out[2]: Q1 Q2 Q3 Q4 Q5 Q6 Q7 Q8 Q9 Q10 ... Q28 Q29 Q30 Q31 Q32 Q33 Q34 Q35 Q36 Cluster 2 ... 117 rows × 37 columns In [4]: | cluster_5 = cluster_5.drop(columns=['Cluster']) cluster 5 Q1 Q2 Q3 Q4 Q5 Q6 Q7 Q8 Q9 Q10 ... Q27 Q28 Q29 Q30 Q31 Q32 Q33 Q34 Q35 Q36 Out[4]: 3 ... 5 5 2 ... 5 117 rows × 36 columns In [5]: # Next the KMO (Kiaser-Meyer-Olkin) test was used to determine whether the cluster is suitable for factor analysis # The closer the overall KMO is to 1, the better suited the dataset. kmo_all, kmo_model = calculate_kmo(cluster_5) print("KMO per variable:", kmo all) print("Overall KMO:", kmo_model) # Since the cluster has a reasonable KMO value, going ahead with the Factor Analysis KMO per variable: [0.8753624 0.71022385 0.93372181 0.85828974 0.85034883 0.86167122 0.88889427 0.91029094 0.89244858 0.86228494 0.63324518 0.60895361 0.90776095 0.82460758 0.79200947 0.9073914 0.9015926 0.895892540.83896281 0.7610136 0.91539386 0.88522506 0.88711132 0.82239618 0.7937945 0.80342156 0.8742284 0.90747866 0.8272541 0.8778996 0.88076142 0.89549342 0.85383294 0.90660564 0.80319907 0.82362279] Overall KMO: 0.8622989636604216 In [6]: # The original correlation matrix for the third cluster: cluster 5.corr() Out[6]: Q 1 Q 2 Q 3 Q 4 Q 5 Q 6 Q 7 Q 8 Q 9 Q 10 ... Q 27 Q 28 Q 29 Q 30 Q 31 Q 32 Q 33 1.000000 0.549596 0.455565 0.430469 0.661759 0.204717 -0.099678 -0.200118 0.580848 -0.069881 ... 0.514901 -0.172580 0.706379 0.181452 -0.160434 0.549596 0.545300 0.228705 0.019986 -0.134238 1.000000 0.317573 0.238204 0.262989 -0.113943 ... 0.205570 -0.129923 0.550601 0.110339 -0.134234 0.266028 -0.047341 -0.0979 0.455565 0.317573 1.000000 0.792832 0.364779 0.252652 0.431952 0.306397 ... 0.505488 0.509330 0.372167 0.524963 0.736146 0.415121 0.255846 0.297 0.238204 0.792832 1.000000 0.474953 0.282873 0.400708 0.758694 0.340266 0.304449 ... 0.374682 0.388240 0.571392 0.422799 0.303 0.545300 0.524963 0.400708 1.000000 0.318334 -0.011559 -0.096733 0.586682 0.031859 ... 0.487347 -0.069026 0.756054 0.181720 -0.052887 0.539596 0.195562 -0.057 0.204717 0.228705 0.736146 0.758694 0.318334 1.000000 0.550445 0.376906 0.242402 0.506419 ... 0.320262 0.396051 0.278704 0.460704 0.553147 0.413676 0.413179 0.442 0.019986 0.364779 0.474953 -0.011559 0.550445 1.000000 0.630134 0.436633 ... -0.002610 0.573789 -0.186565 0.473245 0.619036 0.021620 0.630134 1.000000 -0.007126 -0.200118 -0.134238 0.252652 0.282873 -0.096733 0.376906 0.605115 ... -0.012610 0.728812 -0.181167 0.520657 0.666529 0.118412 0.419423 0.6686 0.262989 0.431952 0.340266 0.021620 -0.007126 0.147352 ... 0.504755 0.586682 0.242402 1.000000 0.037797 0.539855 0.284533 0.094327 0.526853 0.393403 -0.004 -0.113943 0.306397 0.304449 0.436633 0.605115 0.007247 0.357953 0.031859 0.506419 0.147352 1.000000 ... -0.015188 0.579217 0.638239 0.135948 0.5270 0.276292 0.072018 0.124567 0.214306 0.025936 0.240867 0.282218 0.221657 0.166466 ... 0.146010 0.202029 0.165838 0.129177 0.105890 0.156099 0.1282 0.021801 0.001922 0.095299 0.064396 0.078850 0.173455 0.375477 ... 0.159923 0.034803 0.206357 0.145391 0.021329 0.138509 0.001852 0.120610 0.088896 -0.042631 0.103(12 0.486696 0.321482 0.467029 0.404974 0.197499 0.229283 ... 0.386758 0.539106 0.443985 0.046384 0.446520 0.071273 0.413056 0.285125 0.221324 0.353399 0.1673 0.468119 -0.079610 0.228510 -0.058132 0.506464 0.206954 0.004110 0.392438 0.081625 ... -0.037176 0.306078 0.162282 -0.0134 0.029757 0.054173 0.315195 0.295027 0.369771 0.459505 0.278613 0.577794 ... 0.075582 0.360261 0.193011 0.285436 0.467437 0.179769 0.154048 0.365463 0.543926 0.082754 0.488300 ... 0.150150 0.549876 -0.132640 0.450386 0.606632 0.003144 -0.088390 0.319888 0.385959 -0.027609 0.532209 0.569139 0.163727 0.503300 ... 0.202839 0.553500 -0.102112 0.383598 0.591855 0.200058 -0.114469 -0.096060 0.516444 -0.080272 ... 0.606326 0.312730 0.045767 0.497703 0.218258 0.036729 -0.047361 0.359785 0.054143 ... 0.276639 -0.094222 -0.072628 0.390386 0.608079 0.708364 0.146457 0.629588 ... -0.020519 0.748766 -0.091719 0.628581 0.429232 -0.007304 -0.089384 0.582652 0.097442 ... 0.583716 -0.043928 0.640141 0.227787 -0.024093 0.398119 0.508295 0.125208 0.421687 ... -0.043879 0.572043 -0.133212 0.453575 -0.067512 0.269165 0.253294 -0.089518 0.389190 0.420389 0.328246 0.184364 0.271735 ... 0.252759 0.284611 0.398337 0.477140 -0.035665 0.398396 ... -0.186422 0.536759 0.022293 0.272588 0.399203 0.496278 0.057570 0.431354 ... -0.174604 0.548810 -0.164557 0.287890 0.487347 0.320262 -0.002610 -0.012610 0.504755 -0.015188 ... 1.000000 -0.052087 0.582642 0.306043 0.573789 0.728812 0.037797 0.579217 ... -0.052087 1.000000 -0.181606 0.570820 1.000000 -0.186565 -0.181167 0.539855 0.007247 ... 0.582642 -0.181606 0.195845 0.473245 0.520657 0.284533 0.357953 ... 0.306043 0.570820 0.195845 1.000000 0.618772 -0.052887 0.553147 0.619036 0.666529 0.094327 0.638239 ... 0.042470 0.815864 -0.098875 0.618772 0.188606 0.548835 0.076680 0.118412 0.526853 0.135948 ... 0.558214 0.421652 0.419423 0.393403 0.312844 0.495023 0.527677 ... -0.030077 -0.057710 0.442916 0.571728 0.668646 -0.004165 0.724151 -0.142238 0.521452 0.732762 0.405021 0.382746 0.164994 0.501885 0.467918 0.256214 0.418414 0.297134 0.282779 0.416696 0.369089 0.445986 0.360979 0.394537 ... 36 rows × 36 columns In [8]: # The eigenvalues are calculated for the third cluster fa_1 = FactorAnalyzer(n_factors=3, rotation='oblimin').fit(cluster_5) fa_1.loadings_ eigenvalues, _ = fa_1.get_eigenvalues() eigen values = list(eigenvalues) data_ev = pd.DataFrame(eigen_values,columns=['Eigenvalues']) data_ev # In this case, since the first 7 eigenvalues are greater than 1, the no. of factors is take to be 7. **Eigenvalues** Out[8]: 11.622003 7.266139 1.990538 1.758093 1.429721 1.245185 1.065921 0.996036 0.877576 0.711040 10 0.622060 11 0.573227 12 0.535024 0.506471 14 0.461963 0.410649 15 16 0.389377 17 0.372894 18 0.332059 19 0.323559 0.287605 0.258200 22 0.233161 23 0.225169 0.215713 24 25 0.178115 26 0.164256 27 0.162852 28 0.148412 29 0.119574 30 0.113307 31 0.105637 32 0.089093 33 0.078384 34 0.075462 35 0.055524 In [9]: # The factor loadings for the 7 Factors are presented in the form of a heatmap: fa_2 = FactorAnalyzer(n_factors=7, rotation='oblimin').fit(cluster_5) loadings = fa_2.loadings_ plt.figure(figsize=(12, 12)) sns.heatmap(loadings, annot=True, cmap='coolwarm', fmt=".2f") plt.title('Heatmap of Factor Loadings') plt.xlabel('Factors') plt.ylabel('Variables') plt.show() **Heatmap of Factor Loadings** 0.56 -0.23 0.10 0.29 -0.12 0.09 -0.06 0 -0.18 - 0.8 0.66 0.16 -0.08 -0.05 -0.04 0.22 -0.01 0.70 0.19 0.04 -0.02 -0.05 2 -0.80 0.07 0.06 0.09 -0.07 -0.05 -0.04 m --0.09 0.13 -0.04 -0.01 0.80 0.01 0.05 4 --0.02 0.03 0.85 -0.03 -0.02 0.12 0.08 Ω--0.01 0.36 -0.21 0.14 0.24 0.03 0.42 9 -- 0.6 -0.09 0.70 0.04 -0.05 0.08 0.08 0.20 _ -0.13 0.33 -0.19 0.15 0.49 0.05 0.14 ω -0.17 -0.14 0.36 0.15 0.11 -0.02 0.45 6 --0.24 0.39 0.22 0.02 0.15 -0.06 0.33 0. --0.01 -0.17 0.12 0.81 0.05 -0.05 -0.05 - 0.4 -0.12 0.15 0.04 0.42 12 0.51 0.03 0.06 0.79 -0.02 0.07 -0.11 0.18 -0.01 0.05 -0.07 14 -0.35 -0.06 0.13 0.09 -0.01 0.63 0.19 0.41 0.47 -0.08 -0.04 0.17 -0.03 -0.11 0.27 0.29 -0.00 -0.01 0.42 0.04 16 Variables -0.16 0.23 0.16 0.06 -0.03 0.21 - 0.2 18 0.24 0.12 -0.31 0.62 -0.14 0.08 -0.16 19 0.08 -0.12 0.14 0.03 -0.11 0.62 -0.02 20 0.13 0.09 0.23 -0.07 0.11 0.05 0.68 0.24 0.10 -0.48 0.08 0.15 0.39 -0.00 0.19 0.01 0.10 0.45 22 0.12 0.51 -0.03 - 0.0 -0.06 0.06 0.02 0.02 0.10 -0.05 0.69 24 0.13 -0.00 0.18 0.69 0.03 0.09 0.04 25 0.04 0.10 0.19 0.00 0.20 0.70 0.04 26 -0.07 0.04 0.36 -0.45 0.27 0.06 0.22 -0.01 27 -0.04 0.83 0.01 0.03 0.10 0.03 - -0.2 0.17 0.26 -0.21 -0.18 0.11 ω. 0.55 -0.13 29 -0.10 0.62 0.10 0.14 0.26 0.05 -0.15 0.08 0.06 0.18 გ -0.07 -0.13 0.64 0.16 0.23 0.15 0.22 -0.28 -0.01 0.46 -0.10 0.02 0.43 0.05 0.46 -0.09 0.18 -0.16 0.42 0.09 0.13 0.29 0.24 0.03 -0.10 - -0.4 0.15 34 --0.04 0.07 0.08 0.81 -0.04 0.10 -0.02 0.17 0.75 0.17 0.10 -0.01 35 -0.03 0 2 3 5 6 1 4 Factors In [10]: corr_mat_factors_5 = pd.DataFrame(fa_2.phi_) corr_mat_factors_5 # The correlation matrix between the 7 Factors themselves:

2

Out[10]:

0

1 -0.074657

2 0.216908

0.301032

0.402661

In [11]: # Exporting to Excel

0 1.000000 -0.074657 0.216908

0.410532 -0.020930 0.178342

1.000000 0.090221

0.090221 1.000000

0.340310 0.223915

3 0.377905 -0.255720 0.104705 1.000000 -0.061454

0.393148 0.078038 -0.061454

corr_mat_factors_5.to_excel('FA5_KMedoids.xlsx')

3

0.301032

0.393148

0.078038

1.000000

0.256808

0.423330

0.377905

-0.255720

0.104705

0.180937

0.059020

6

5

0.410532 0.402661

-0.020930 0.340310

0.178342 0.223915

0.256808 0.423330

1.000000 0.379578

0.379578 1.000000