OPERATION RESEARCH

The criteria for the pre-selection are:

• Viscosity of the neat resin should be less than 1 Pa·s (1000 cps) to enable VARIM processing. • Tg (glass transition temperature) of the neat resin should preferable be higher than 50-60°C. • The process temperature should be lower than 230°C, so that low cost accessories can be used. • The resin should have a long pot life. • Cost of the resin should be affordable. • Availability of basic knowledge and technology.

Algorithm

- 1. Elimination Search according to constraints d
- 2. Obtain the decision matrix and relative importance matrix a
- 3. Normalize the decision matrix using the method from TOPSIS rij=xij/Σ(xij)^0.5
- 4. Normalize the relative importance matrix using the Geometric Mean from AHP Method GMi=(πaij)^(1/n) by using AHP
- 5. Eigen value calculation and multiplying decision matrix by weights to obtain normalized weighted matrix
- 6. Obtain the best and worst solution from weighted normalized matrix using TOPSIS
- 7. Obtain best soln using calculating the Euclidean distance and then ranking them according to RSI score.RSI=S_minus/S_plus+S_minus

In [1]:

```
#Using the suitable libraries
import pandas as pd
import numpy as np
import matplotlib.pyplot as plt
import plotly.graph_objects as go
%matplotlib inline
```

In [2]:

```
#Loading the Excel sheet containing the Decision matrix
df=pd.read_excel("")
df
```

Out[2]:

	Attributes Material	Viscosity (Pa s)	Processing temperature(Celsius)\n	Cost(€/kg)	Availability
0	Polymethylmethacrylate (PMMA)	0.10	140	4.0	0.75
1	Polyamide 6 (PA 6)	0.01	150	2.5	1.00
2	Polyamide 12 (PA 12)	0.05	220	20.0	0.75
3	Polyethylene Terephtalate (PET)	0.95	280	3.0	0.50
4	Thermoplastic Polyurethane (TPU)	0.80	300	10.0	0.75
5	Polybutylene Terephtalate (PBT)	0.02	200	9.0	1.00
6	Polycarbonate (PC)	0.70	250	5.0	0.25
7	Polyether Ketone (PEK)	0.10	350	60.0	0.25
8	Polyphthalamide (PPA)	1.00	250	5.5	0.75

```
In [3]:
```

```
d=df.drop(columns=df.columns[0])
```

In [4]:

Out[4]:

```
Viscosity (Pa s) Processing temperature(Celsius)\n Cost(€/kg) Availability
0
              0.10
                                                    140
                                                                            0.75
                                                                 4.0
1
              0.01
                                                    150
                                                                 2.5
                                                                             1.00
              0.05
2
                                                    220
                                                                20.0
                                                                            0.75
3
              0.95
                                                    280
                                                                 3.0
                                                                            0.50
              0.80
                                                    300
                                                                            0.75
4
                                                                10.0
              0.02
5
                                                    200
                                                                 9.0
                                                                            1.00
              0.70
                                                    250
                                                                            0.25
6
                                                                 5.0
7
              0.10
                                                    350
                                                                60.0
                                                                            0.25
              1.00
                                                    250
8
                                                                 5.5
                                                                            0.75
```

In [5]:

```
#Normalizing the decision matrix
def den(i,j,d):
    sum_=0
    for i in range(0,d.shape[0]):
        sum_+=(d[i][j])**2
    return sum_
d=np.array(d)
r=np.zeros((d.shape[0],d.shape[1]))
for i in range(0,d.shape[0]):
    for j in range(0,d.shape[1]):
        r[i][j]=d[i][j]/(den(i,j,d))**0.5
```

In [6]:

r

Out[6]:

```
array([[0.05720828, 0.18939682, 0.06123126, 0.34874292], [0.00572083, 0.20292517, 0.03826954, 0.46499055], [0.02860414, 0.29762358, 0.30615632, 0.34874292], [0.54347862, 0.37879365, 0.04592345, 0.23249528], [0.4576662, 0.40585034, 0.15307816, 0.34874292], [0.01144166, 0.27056689, 0.13777034, 0.46499055], [0.40045793, 0.33820861, 0.07653908, 0.11624764], [0.05720828, 0.47349206, 0.91846896, 0.11624764], [0.57208275, 0.33820861, 0.08419299, 0.34874292]])
```

In [7]:

```
#Loading the Excel sheet containing the realtive importance matrix
dfl=pd.read_excel("")
dfl
```

Out[7]:

Unnamed: 0 Viscosity Processing temp cost availability

0	Viscosity	1	1	1	1
1 F	Processing temp	1	1	1	1
2	cost	1	1	1	1
3	availability	1	1	1	1

In [8]:

```
df1.drop(columns=['Unnamed: 0'],inplace=True)
```

```
df1
Out[9]:
  Viscosity Processing temp cost availability
0
                                   1
                          1
                                   1
1
        1
                     1
2
3
        1
                     1
                          1
                                   1
In [10]:
#To remove the bias, incorporating the AHP method for calcualtion of weights
def gm(df1):
    gmi=np.ones((df1.shape[0]))
    for i in range(0,df1.shape[0]):
        for j in range(0,df1.shape[1]):
            gmi[i] *=df1[i][j]
        gmi[i] = gmi[i] * * (1/df1.shape[1])
    return gmi
df1 np=np.array(df1)
gmi=gm(df1 np)
w=np.zeros(gmi.shape[0])
for i in range(0,gmi.shape[0]):
    w[i] = gmi[i] / sum(gmi)
N2=np.transpose(w)
N3=np.dot(df1 np, N2)
N4=N3/N2
N2, N3, N4, gmi, df1 np, w
Out[10]:
(array([0.25, 0.25, 0.25, 0.25]),
 array([1., 1., 1., 1.]),
 array([4., 4., 4., 4.]),
 array([1., 1., 1., 1.]),
 array([[1, 1, 1, 1],
        [1, 1, 1, 1],
        [1, 1, 1, 1],
        [1, 1, 1, 1]], dtype=int64),
 array([0.25, 0.25, 0.25, 0.25]))
In [11]:
#Calculating the eigen values using the relative importance matrix
eigen, righteigen=np.linalg.eig(df1 np)
In [12]:
eigen, righteigen
Out[12]:
(array([0.0000000e+00, 4.00000000e+00, 0.00000000e+00, 2.80731443e-32]),
                                  , -0.8660254 , -0.64641535],
 array([[-0.8660254 , -0.5
        [ 0.28867513, -0.5
                                     0.28867513, -0.32788993],
        [ 0.28867513, -0.5
                                     0.28867513, 0.48715264],
        [ 0.28867513, -0.5
                                  , 0.28867513, 0.48715264]]))
In [13]:
lambdamax=max(eigen)
lambdamax
```

In [9]:

Out[13]:

```
In [14]:
#Calculating the consistency ratio to ensure best solution. CR=CI/RI. CR<=1
#Random index values corresponding to different matrix sizes in array ri
ri=np.array([0.00,0.00,0.58,0.90,1.12,1.24,1.32,1.41,1.45,1.49])
ci=(lambdamax-df1 np.shape[0])/(df1 np.shape[0]-1)
consistency ratio=ci/ri[df1 np.shape[0]-1]
consistency ratio
Out[14]:
0.0
In [15]:
#Obtaining the weighted normalized matrix
V = r * w
\nabla
Out[15]:
array([[0.01430207, 0.04734921, 0.01530782, 0.08718573],
       [0.00143021, 0.05073129, 0.00956739, 0.11624764],
       [0.00715103, 0.07440589, 0.07653908, 0.08718573],
       [0.13586965, 0.09469841, 0.01148086, 0.05812382],
       [0.11441655, 0.10146258, 0.03826954, 0.08718573],
       [0.00286041, 0.06764172, 0.03444259, 0.11624764],
       [0.10011448, 0.08455215, 0.01913477, 0.02906191],
       [0.01430207, 0.11837301, 0.22961724, 0.02906191],
       [0.14302069, 0.08455215, 0.02104825, 0.08718573]])
In [16]:
#Obtaining the best and worst system from matrix V
V plus=np.zeros(V.shape[1])
V minus=np.zeros(V.shape[1])
for j in range(0, V.shape[1]):
    V plus[j]=np.max(V[:,[j]])
    V minus[j]=np.min(V[:,[j]])
V_plus, V_minus
Out[16]:
(array([0.14302069, 0.11837301, 0.22961724, 0.11624764]),
 array([0.00143021, 0.04734921, 0.00956739, 0.02906191]))
In [17]:
#Caluculating the Euclidean distance from best and worst system of each entry in V. Then
Ranking them according to RSI
#RSI=S minus/S plus+S minus
S_plus=np.zeros(V.shape[0])
S minus=np.zeros(V.shape[0])
RSI=np.zeros(V.shape[0])
for i in range(0, V.shape[0]):
    for j in range(0, V.shape[1]):
        S plus[i]+=(V[i][j]-V plus[j])**2
        S minus[i]+=(V[i][j]-V minus[j])**2
    S plus[i]=S plus[i]**0.5
    S minus[i]=S minus[i]**0.5
    RSI[i]=S minus[i]/(S plus[i]+S minus[i])
```

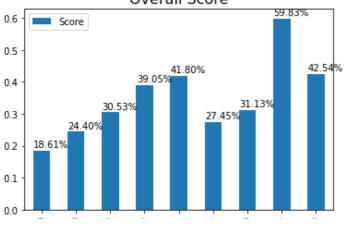
```
In [18]:
```

```
RSI
```

Out[18]:

```
array([0.18613542, 0.24404589, 0.30531042, 0.39046714, 0.41802793, 0.27449964, 0.31129606, 0.5983321 , 0.42544464])
```

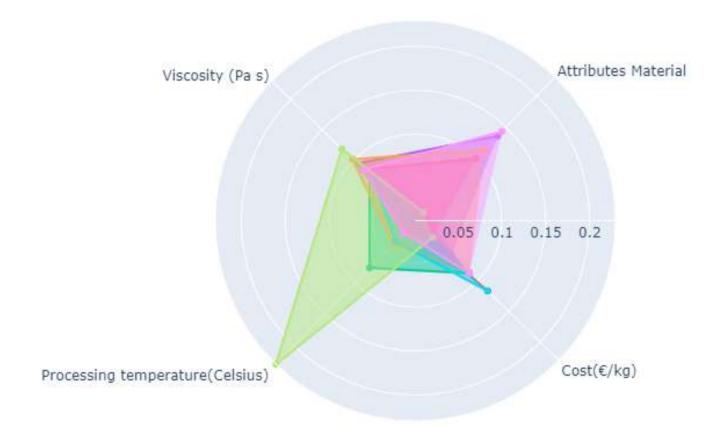
```
In [19]:
#Sorting the items according to the RSI Score
sortedRSI=np.sort(RSI)[::-1]
sortedRSI index=np.zeros(RSI.shape[0])
for i in range(0,RSI.shape[0]):
    sortedRSI index[i]=int(np.where(RSI==sortedRSI[i])[0]+1)
sortedRSI index=sortedRSI index.astype(int)
RSI, sortedRSI, sortedRSI index
Out[19]:
(array([0.18613542, 0.24404589, 0.30531042, 0.39046714, 0.41802793,
        0.27449964, 0.31129606, 0.5983321 , 0.42544464]),
array([0.5983321 , 0.42544464, 0.41802793, 0.39046714, 0.31129606,
        0.30531042, 0.27449964, 0.24404589, 0.18613542]),
array([8, 9, 5, 4, 7, 3, 6, 2, 1]))
In [20]:
#Obtaining the list of items in best to worst order and obtaining the bar graph to show c
omparison b/w different items
names=list(df.columns)
items=list(df[names[0]])
for i in range(0, sortedRSI index.shape[0]):
     print(items[sortedRSI_index[i]-1])
df final = {names[0]:items, 'Score': RSI}
df final= pd.DataFrame(data=df final)
ax=df final.plot(kind='bar')
ax.set title("Overall Score", fontsize=16)
for p in ax.patches:
    ax.annotate("{:.2%}".format(p.get height()),
                xy = (p.get x() + 0.02, p.get height() + 0.01))
ax.set xticklabels(items)
Polyether Ketone (PEK)
Polyphthalamide (PPA)
Thermoplastic Polyurethane (TPU)
Polyethylene Terephtalate (PET)
Polycarbonate (PC)
Polyamide 12 (PA 12)
Polybutylene Terephtalate (PBT)
Polyamide 6 (PA 6)
Polymethylmethacrylate (PMMA)
Out [20]:
[Text(0, 0, 'Polymethylmethacrylate (PMMA)'),
Text(1, 0, 'Polyamide 6 (PA 6)'),
Text(2, 0, 'Polyamide 12 (PA 12)'),
Text(3, 0, 'Polyethylene Terephtalate (PET)'),
Text(4, 0, 'Thermoplastic Polyurethane (TPU)'),
Text(5, 0, 'Polybutylene Terephtalate (PBT)'),
Text(6, 0, 'Polycarbonate (PC)'),
Text(7, 0, 'Polyether Ketone (PEK)'),
Text(8, 0, 'Polyphthalamide (PPA)')]
                 Overall Score
                                    59.83%
0.6
       Score
```



```
Polymethylmethacrylate (PMMA)
Polyamide 12 (PA 6)
Polyethylene Terephtalate (PET)
Thermoplastic Polyurethane (TPU)
Polybutylene Terephtalate (PBT)
Polybutylene Terephtalate (PBT)
Polybutylene Terephtalate (PBT)
Polybutylene (PEK)
```

In [21]:

```
#For obtaining the spider diagram also known as the Radar plot
list name = names
attributes= items
fig = go.Figure()
for i in range(0, V.shape[0]):
    fig.add trace(go.Scatterpolar(
          r=V[i],
          theta=list name,
          fill='toself',
          name=attributes[i]
    ) )
fig.update_layout(
  polar=dict(
    radialaxis=dict(
      visible=True,
      range=[np.min(V), np.max(V)]
    )),
  showlegend=True
fig.show()
```



- Polymethylmethacrylate (PMMA)
- Polyamide 6 (PA 6)
- Polyamide 12 (PA 12)
- Polyethylene Terephtalate (PET)
 - Thermoplastic Polyurethane (TPU)
- Polybutylene Terephtalate (PBT)
- Polycarbonate (PC)
- Polyether Ketone (PEK)
- Polyphthalamide (PPA)