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```
import numpy as np
proto_nodes = np.array([
                        [0,0],
                        [0.5,1],
                        [1,0.5]
1)
def manhattan distance(x,y):
  return np.absolute(x[0] - y[0]) + np.absolute(x[1] - y[1])
def rb function(x, p node):
  return max(0, 1 - manhattan_distance(x, p_node))
datapoint = np.array([0.6,0.8])
o nodes = 3
weights = np.zeros([proto nodes.shape[0], o nodes])
weights[:,0] = 0.6
weights[:,1] = -0.4
weights[:,2] = 0
weights
    array([[ 0.6, -0.4, 0. ],
           [0.6, -0.4, 0.],
           [0.6, -0.4, 0.]
rbf_results = []
for n in proto_nodes:
  rbf results.append(rb function(datapoint, n))
rbf_results = np.array(rbf_results)
outputs = np.dot(rbf_results, weights)
outputs
    array([ 0.6, -0.4, 0. ])
```

Class A is the winner!

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