Q1. PCA

1.

Given the points (i,i), (i, i+1) for i = 1,, 10 Let us represent the given space using a 20 x 2 matrix

1	1
2	2
3	3
4	4
5	5
6	6
7	7
8	8
9	9
10	10
1	2
2	3
3	4
4	5
5	6
6	7
7	8
8	9
9	10
10	11

Mean of x-coordinates

$$x_{mean} = (1 + 2 + ... + 10 + 1 + 2 + + 10) / 20$$

= 5.5

Mean of y-coordinates

$$y_{mean} = (1 + 2 + ... + 10 + 2 + 3 + + 11) / 20$$

= 6

Now, for computing the centered data points, we will shift all points using

$$xj_{centered} = xj - x_{mean}$$

 $yj_{centered} = yj - x_{mean}$

where, (xj, yj) are the j-th points in the given space.

Using above, the $X_{centered}$ 20 x 2 matrix can be represented as follow:

2. S =
$$(1/n-1)X^{T}X$$

For our centered data, $X_{\it centered}$ as computed in (1) can be used.

n = 20

(not displaying the matrix multiplication because of large dimensionality)

Upon computation,

cov =

```
(8.684210526315790 -8.684210526315790)
8.684210526315790 8.947368421052632)
```

3. Using eigen computation, eigenvalues and eigenvectors of cov matrix are

```
λ1 = 17.500996753004870

v1 = (0.701730084756145, 0.712442901675730)

λ2 = 0.130582194363553

v2 = (-0.712442901675730, 0.701730084756145)
```

Now, projection of a point X on a vector v is given by, $X_{centered}v^T$

Thus, projection of any given points on these two vectors would be $(X_{centered}v_1^{\ T},\ X_{centered}v_2^{\ T})$

Using this for all the centered points, we can compute the projected values to be the following 20 x 2 matrix. (next page)

original (x,y)	projected x	projected y
(1,1)	-6.71999989	-0.3026573662
(2,2)	-5.305826903	-0.3133701832
(3,3)	-3.891653917	-0.3240830001
(4,4)	-2.47748093	-0.334795817
(5,5)	-1.063307944	-0.3455086339
(6,6)	0.3508650424	-0.3562214508
(7,7)	1.765038029	-0.3669342678
(8,8)	3.179211015	-0.3776470847
(9,9)	4.593384002	-0.3883599016
(10,10)	6.007556988	-0.3990727185
(1,2)	-6.007556988	0.3990727185
(2,3)	-4.593384002	0.3883599016
(3,4)	-3.179211015	0.3776470847
(4,5)	-1.765038029	0.3669342678
(5,6)	-0.3508650424	0.3562214508
(6,7)	1.063307944	0.3455086339
(7,8)	2.47748093	0.334795817
(8,9)	3.891653917	0.3240830001
(9,10)	5.305826903	0.3133701832
(10,11)	6.71999989	0.3026573662

Q2. Resolution

```
Knowledge Base: p \Rightarrow (q \Rightarrow r) \dots (0)
Representing knowledge base in CNF, we get
p \Rightarrow (\neg q \lor r) \dots using implication elimination
¬p v (¬q v r) ...... using implication elimination
(¬p v ¬q v r) ..... using associativity of v
To Prove: (p \land q) => (q => r) \dots (1)
So, we can prove that (0) entails (1), i.e.,
Whenever (0) is true (which is given), then (1) is also true.
Representing (1) in CNF, we get
(p \land q) => (\neg q \lor r) \dots using implication elimination
(\neg(p \land q)) \lor (\neg q \lor r) ..... using implication elimination
(¬p v ¬q) v (¬q v r) ...... using de Morgan
(¬p v ¬q v ¬q v r) ...... using associativity of v
(\neg p \lor \neg q \lor r)
Now, negation of (1) would be,
¬(¬p v ¬q v r)
(p \land q \land \neg r) ...... using de Morgan
Adding negation of (1) to Knowledge Base, we get
a1: ¬p v ¬q v r
b1: p
b2: q
b3: ¬r
Goal: empty set
Step 1: resolve a1, b1:
                                ¬q∨r
Step 2: resolve above, b2:
Step 3: resolve above, b3:
                                empty
Thus, (0) entails (1)
Hence proved.
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Q3. Hierarchical Clustering

1. Calculating the euclidean distance between all the points, $d = ((x1-x2)^2 + (y1-y2)^2)^{1/2}$

	Madison (-89, 43)	Seattle (-122, 48)	Boston (-71, 42)	Vancouver (-123, 49)	Winnipeg (-97, 50)	Montreal (-74, 46)
Madison (-89, 43)	0	33.376639	18.027756	34.525353	10.630146	15.297059
Seattle (-122, 48)	33.376639	0	51.351728	1.414214	25.079872	48.041649
Boston (-71, 42)	18.027756	51.351728	0	52.469038	27.202941	5
Vancouver (-123, 49)	34.525353	1.414214	52.469038	0	26.019224	49.091751
Winnipeg (-97, 50)	10.630146	25.079872	27.202941	26.019224	0	23.345235
Montreal (-74, 46)	15.297059	48.041649	5	49.091751	23.345235	0

Iteration 1:

Closest pair of clusters => (Vancouver, Seattle)

Distance between them as defined by complete linkage => 1.414214

All clusters at the end of that iteration =>

(Vancouver, Seattle), (Madison), (Boston), (Winnipeg), (Montreal)

Iteration 2:

	Madison (-89, 43)	Boston (-71, 42)	Winnipeg (-97, 50)	Montreal (-74, 46)	(Vancouver, Seattle)
Madison (-89, 43)	0	18.027756	10.630146	15.297059	34.525353
Boston (-71, 42)	18.027756	0	27.202941	5	52.469038
Winnipeg (-97, 50)	10.630146	27.202941	0	23.345235	26.019224
Montreal (-74, 46)	15.297059	5	23.345235	0	49.091751

Closest pair of clusters => (Montreal, Boston)
Distance between them as defined by complete linkage => 5
All clusters at the end of that iteration => (Vancouver, Seattle), (Montreal, Boston), (Madison), (Winnipeg)

Iteration 3:

	Madison (-89, 43)	Winnipeg (-97, 50)	(Vancouver, Seattle)	(Montreal, Boston)
Madison (-89, 43)	0	10.630146	34.525353	18.027756
Winnipeg (-97, 50)	10.630146	0	26.019224	27.202941
(Montreal, Boston)	18.027756	27.202941	52.469038	0

Closest pair of clusters => (Madison, Winnipeg)
Distance between them as defined by complete linkage => 10.630146
All clusters at the end of that iteration => (Vancouver, Seattle), (Montreal, Boston), (Madison, Winnipeg)

Iteration 4:

	(Vancouver, Seattle)	(Montreal, Boston)	(Madison, Winnipeg)
(Montreal, Boston)	52.469038	0	27.202941
(Vancouver, Seattle)	0	52.469038	34.525353
(Madison, Winnipeg)	34.525353	27.202941	0

Closest pair of clusters => ((Madison, Winnipeg), (Montreal, Boston))
Distance between them as defined by complete linkage => 27.202941
All clusters at the end of that iteration => (Vancouver, Seattle), (Madison, Winnipeg, Montreal, Boston)

2. Calculating the euclidean distance between all the points, constraint being US to Canada city distance = ND (infinite),= $d = ((x1-x2)^2 + (y1 - y2)^2)^{1/2}$

$$d = ((x1 - x2)^2 + (y1 - y2)^2)^{1/2}$$

	Madison (-89, 43)	Seattle (-122, 48)	Boston (-71, 42)	Vancouver (-123, 49)	Winnipeg (-97, 50)	Montreal (-74, 46)
Madison (-89, 43)	0	33.376639	18.027756	ND	ND	ND
Seattle (-122, 48)	33.376639	0	51.351728	ND	ND	ND
Boston (-71, 42)	18.027756	51.351728	0	ND	ND	ND
Vancouver (-123, 49)	ND	ND	ND	0	26.019224	49.091751
Winnipeg (-97, 50)	ND	ND	ND	26.019224	0	23.345235
Montreal (-74, 46)	ND	ND	ND	49.091751	23.345235	0

Iteration 1:

Closest pair of clusters => (Madison, Boston)

Distance between them as defined by complete linkage => 18.027756

All clusters at the end of that iteration => (Madison, Boston), Seattle, Vancouver, Winnipeg, Montreal

Iteration 2:

	Seattle (-122, 48)	Vancouver (-123, 49)	Winnipeg (-97, 50)	Montreal (-74, 46)	(Madison, Boston)
Seattle (-122, 48)	0	ND	ND	ND	51.351728
Vancouver (-123, 49)	ND	0	26.019224	49.091751	ND
Winnipeg (-97, 50)	ND	26.019224	0	23.345235	ND
Montreal (-74, 46)	ND	49.091751	23.345235	0	ND

Closest pair of clusters => (Winnipeg, Montreal)
Distance between them as defined by complete linkage => 23.345235
All clusters at the end of that iteration => (Madison, Boston), (Winnipeg, Montreal), Seattle, Vancouver

Iteration 3:

	Seattle (-122, 48)	Vancouver (-123, 49)	(Madison, Boston)	(Winnipeg, Montreal)
Seattle (-122, 48)	0	ND	51.351728	ND
Vancouver (-123, 49)	ND	0	ND	49.091751
(Madison, Boston)	51.351728	ND	0	ND

Closest pair of clusters => (Vancouver, Winnipeg, Montreal)
Distance between them as defined by complete linkage => 49.091751
All clusters at the end of that iteration => (Madison, Boston), (Vancouver, Winnipeg, Montreal), Seattle

Iteration 4:

	Seattle (-122, 48)	(Madison, Boston)	(Vancouver, Winnipeg, Montreal)
Seattle (-122, 48)	0	51.351728	ND
(Madison, Boston)	51.351728	0	ND
(Vancouver, Winnipeg, Montreal)	ND	ND	0

Closest pair of clusters => (Madison, Boston, Seattle)
Distance between them as defined by complete linkage => 51.351728
All clusters at the end of that iteration => (Madison, Boston, Seattle), (Vancouver, Winnipeg, Montreal)

Q4. K-means Clustering

1.

Iteration 1:

	dist_c1 (0)	dist_c2 (9)	у
x1 (10)	10	1	y1 = 2
x2 (8)	8	1	y2 = 2
x3 (6)	6	3	y3 = 2
x4 (4)	4	5	y4 = 1
x5 (3)	3	6	y5 = 1
x6 (2)	2	7	y6 = 1

Updated centers

$$c1 = (x4 + x5 + x6) / 3$$
$$= (4 + 3 + 2) / 3$$
$$= (9) / 3$$
$$= 3$$

$$c2 = (x1 + x2 + x3) / 3$$

$$= (10 + 8 + 6) / 3$$

$$= (24) / 3$$

$$= 8$$

Energy =
$$d(x1, c2)^2 + d(x2, c2)^2 + d(x3, c2)^2 + d(x4, c1)^2 + d(x5, c1)^2 + d(x6, c1)^2$$

= $1^2 + 1^2 + 3^2 + 4^2 + 3^2 + 2^2$
= $1 + 1 + 9 + 16 + 9 + 4$
= 40

Iteration 2:

	dist_c1 (3)	dist_c2 (8)	у
x1 (10)	7	2	y1 = 2
x2 (8)	5	0	y2 = 2
x3 (6)	3	2	y3 = 2
x4 (4)	1	4	y4 = 1
x5 (3)	0	5	y5 = 1
x6 (2)	1	6	y6 = 1

As clustering remains the same as Iteration 1, implying the centroids have converged, hence we stop here.

Energy =
$$d(x1, c2)^2 + d(x2, c2)^2 + d(x3, c2)^2 + d(x4, c1)^2 + d(x5, c1)^2 + d(x6, c1)^2$$

= $2^2 + 0^2 + 2^2 + 1^2 + 0^2 + 1^2$
= $4 + 0 + 4 + 1 + 0 + 1$
= 10

2.

Iteration 1:

	dist_c1 (8)	dist_c2 (9)	у
x1 (10)	2	1	y1 = 2
x2 (8)	0	1	y2 = 1
x3 (6)	2	3	y3 = 1
x4 (4)	4	5	y4 = 1
x5 (3)	5	6	y5 = 1
x6 (2)	6	7	y6 = 1

Updated centers

$$c1 = (x2 + x3 + x4 + x5 + x6) / 5$$

$$= (8 + 6 + 4 + 3 + 2) / 5$$

$$= (23) / 5$$

$$= 4.6$$

$$c2 = (x1) / 1$$

= (10) / 1
= 10

Energy =
$$d(x1, c2)^2 + d(x2, c2)^2 + d(x3, c2)^2 + d(x4, c1)^2 + d(x5, c1)^2 + d(x6, c1)^2$$

= $1^2 + 0^2 + 2^2 + 4^2 + 5^2 + 6^2$
= $1 + 0 + 4 + 16 + 25 + 36$
= 82

Iteration 2:

	dist_c1 (4.6)	dist_c2 (10)	у
x1 (10)	5.4	0	y1 = 2
x2 (8)	3.4	2	y2 = 2
x3 (6)	1.4	4	y3 = 1
x4 (4)	0.6	6	y4 = 1
x5 (3)	1.6	7	y5 = 1
x6 (2)	2.6	8	y6 = 1

Updated centers

$$c1 = (x3 + x4 + x5 + x6) / 4$$

$$= (6 + 4 + 3 + 2) / 4$$

$$= (15) / 4$$

$$= 3.75$$

$$c2 = (x1 + x2) / 2$$

= $(10 + 8) / 2$
= $(18) / 2$
= 9

Energy =
$$d(x1, c2)^2 + d(x2, c2)^2 + d(x3, c2)^2 + d(x4, c1)^2 + d(x5, c1)^2 + d(x6, c1)^2$$

= $0^2 + 2^2 + 1.4^2 + 0.6^2 + 1.6^2 + 2.6^2$
= $0 + 4 + 1.96 + 0.36 + 2.56 + 6.76$
= 15.64

Iteration 3:

	dist_c1 (3.75)	dist_c2 (9)	у
x1 (10)	6.25	1	y1 = 2
x2 (8)	4.25	1	y2 = 2
x3 (6)	2.25	3	y3 = 1
x4 (4)	0.25	5	y4 = 1
x5 (3)	0.75	6	y5 = 1
x6 (2)	1.75	7	y6 = 1

As clustering remains the same as Iteration 2, implying the centroids have converged, hence we stop here.

Energy =
$$d(x1, c2)^2 + d(x2, c2)^2 + d(x3, c2)^2 + d(x4, c1)^2 + d(x5, c1)^2 + d(x6, c1)^2$$

= $1^2 + 1^2 + 2.25^2 + 0.25^2 + 0.75^2 + 1.75^2$
= $1 + 1 + 5.0625 + 0.0625 + 0.5625 + 3.0625$
= 10.75

3.

As the final energy in part 1 is less, hence starting with centers c1 (0) and c2 (9) is providing a better minima of distortion. Thus the k-means solution of part 1 is better.