PRML Mid Semester Solution

Note: Marking Scheme is at the end.

1(a)

(i)

Surface plot of $f_1(x,y) = x_1^2 + x_2^2 + 2x_1 + 2x_2$

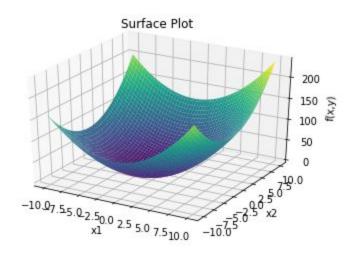


Figure: 1

It is a convex surface.

Surface plot of $f_2(x,y) = x_1 sin(x_1) + x_2 sin(x_2)$

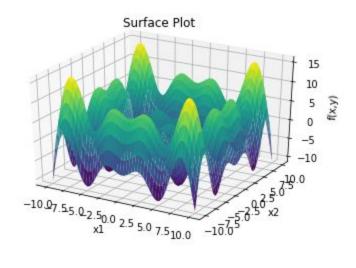


Figure: 2

It is a non-convex surface.

(iv) For $f_1(x,y)$

Initialization Learning rate		location of min	Comment				
-9,-9	0.001	-1.00, -1.00					
-5,-5	0.001	-1.00, -1.00	Late convergence compared to I.r =0.1				
7,-8	0.001 -1.00, -1.00						
-9,-9	0.1	-1.00, -1.00					
-5,-5	0.1 -1.00, -1.00		Converge to global min irrespective of initialization.				
7,-8	0.1	-0.99, -1.00					
-9,-9	1.1	-1.21e+80, -1.21e+80					
-5,-5	1.1 -6.07e+79, -6.07e+79		Does not converge when I.r. is high.				
7,-8	1.1	1.21e+80, -1.06e+80	11 00				

For $f_2(x,y)$

Initialization Learning rate		location of min	Comment			
-1,-1 0.001		0,0				
-5,-5	0.001	-4.91, -4.91	Late convergence as the I.r. is low. Difference poin of convergence when initialization point is different			
7,-8	0.001	4.92, -10	or convergence when militalization point is different.			
-1,-1	0.1	0,0	Difference point of convergence when initialization			
-5,-5	0.1	-4.91, -4.91	point is different. This is because the gradient			
7,-8	0.1	4.91, -10	decent algorithm gets stuck at local minima.			
-1,-1	2.5	7.11e+107 7.11e+107				
-5,-5	2.5	1.82e+105 1.82e+105	Does not converge when l.r is high.			
7,-8	2.5	-1.04e+120 -5.92e+107	0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0			

From Figure 1 we can see that their exists (ii)a minima here: 4,(x1)x2) = x12 + x2 +2x1+2x2. For binding location of minima:- $\frac{\partial f_i(x_1, x_2)}{\partial x_i} = 2x_1 + 2 = 0, \quad \frac{\partial f_i(x_1, x_2)}{\partial x_i} = 2x_2 + 2 = 0$

=) $\chi_1 = -1$, $\chi_2 = -1$

Mence, (x, 1x,) = (-1, -1) is the Location of global total minima.

From Figure 2 it is clear that their exists multiple local maxima e local numa minima.

Mence, for binding minim location of minimum volue use need to bind value at all possible local minima's l'extremes values.

fr (2, 12) = M, sinx, + M2 ein M2 21/2

2 (γι, χ) = χ, (ω) χ, + sinχ, βχ (γι, χ) = χ, (ω) χ, + sinχ, χ ∈ [-10,10].

(1) Roll of Contingence Chan classing said is

English Man Jacobs Comment of the

 $\frac{\partial f_{\mathbf{Q}}(\mathbf{x}_{1}, \mathbf{x}_{0})}{\partial \mathbf{x}_{1}} = 0 = \mathbf{x}_{1}, \quad \cos \mathbf{x}_{1} + \sin \mathbf{x}_{1}$ $\Rightarrow \mathbf{x}_{1} = -\tan \mathbf{x}_{1}$ $\Rightarrow \mathbf{x}_{1} = 0, \pm 7, \pm 25, \pm 4, 493$

 $\frac{\partial f_2(x_1, x_2)}{\partial x_2} = x_2 \cos x_2 + x_2 = 0$

=> 1/2 = -ton/2

=) x3 = 0, ± 4.493, ± 7.725.

Therefore, there are 25 points where local manima or local minima tom exists.

We also need to check at the f 4 extremas (-10, -10), (-10, 10), (+10, -10), (105,10).

On checking on all 29 points use get the minimum at extremum ; e (\$10, ±10).

(iii) horadient Descent

Pros

(a) Always converges to global minima when surface is conven ind learning state is small.

(b) Rate of convergence depends whon the charice of learning rate:

Cons

When the Surface is non-conven then reaching toid optiobal min defends whom intealigation.

When learning rate is large then there is divergence. When I. or is long time to winverge to the min.

Normal Equation

Pros

(a) Inloses to for small deterset

(b) hives parameters values in one step

Spired Spirel

LEW M. MUS SAR

Colory Es:

Cons

Fails for largers dataset.

For parameter computation in one state, inverse of nxn matrix is computed and this is computationally expensive. Regoldsion

(i) 9t products continuous crowbs the data according values and their to the similar data points.

(ii) It is superiorised learning technique learning technique

(iii) Eg:- Predictory Predicting which point person's income based on group cluster new data various features

Regression

(cost for: - 1 2 (ho(xi) - yi))2

ho(x) - hypothesis for

y - output:

M - # Examples.

As the Surface is convon it weill reach the minima provided ho(2) is a linear function.

Cost fn: 1 \(\subsetence \in \) \(\lambda \) \(\lambda

The second summation in the above equation is like identity function and it is one only when the point; lectoraps to the ith cluster. This identity brings non linearity in cost for plance, it is a non-wowen for.

户

Performing clustering task for

Data = 0.5, 0.8, 0.9, 1.0, 1.1, 1.2

Let's take 0.8 and 0.9 as cluster centers

Philial centers > ... $\mu_{i}^{(i)} = 0.8$, $\mu_{i}^{(i)} = 0.9$

Calculating distance

vacco		}	- 4	-	[1,1]	1.2
Data	0.5	0.8	0.9	1.0		
(4°) 0.8		0	0-1	0.2	0-3	0.4
(2)	0.4	0-1	0	0-1	0.2	0.31
1, 30-9						

Selecting points to the clusters which are at

(0.9, 1.0, 1.1, 1.2)

New cluster center

$$\mu_1^{\text{new}} = 0.5 + 0.8 = 0.65$$
, $\mu_2^{\text{new}} = 0.9 + 1.1 + 1.11$

$$= (0.5-0.65)^{2} + (0.8-0.65)^{2} + (0.9-1.05)^{2} + (1.0-1.05)^{2} + (1.2-1.05)^{2} + (1.2-1.05)^{2} + (1.2-1.05)^{2} + (1.2-1.05)^{2}$$

$$= 0.095 - 0$$

Distance Metric Wisit ginew, ynew

New : Cluster center => @

Stopping coileria >

As the mean do not change we will Stop at this step.

Checkering using initial centers as $u_1^{(i)} = 0.5$, $u_2^{(i)} = 1.0$.

N	$u_{i}^{(c)}$ =	0.5	, ,	(i) =	1.0.	(1011/11
Data 4(1)=0.5	6.5:	0.8	D.9	1.0	1:1-2	1:2
11(1)	В	0.3	0.4	0.5	0.6	0.7
11(i) = 1.0	0.5	0.5	0.1	0	b.1	0.2

(0.5)

12 (0-8, 0-9, 1-0, 1-1, 1-2)

New Cluster centers:- $U_1^{(new)} = 0.8 + 0.9 + 1.0 + 1.1 + 1.2 = 1.0$

 $(ext = (0.5-0.5)^{2} + (0.8-1.0)^{2} + (0.9-1.0)^{2} + (1.0-1.0)^{2} + (1.1-1.0)^{2} + (1.2-1.0)^{2}$ $(1.1-1.0)^{2} + (1.2-1.0)^{2}$ $(1.1-1.0)^{2} + (1.2-1.0)^{2}$

The cluster center ded not change hence, were stop at this stop.

Cluster 1 + 0.5, Cluster 2 - (0.6, 0.9, 1, 1.1, 1.7)

We can observe that with different initialization we get different cluster.

Mence the for is not a wonven function & reaches different local minima when initialized differently.

So, for landing the lost cluster we do multiple random initialization & bind the cluster with minimum last (Error.

Here In this example dusters with initial centers.

(i) from the t-SME plot of the balanced dataset we see that two proper dusters are formed

Checause when compared to GMM lecause
the circular clusters well easily easily
cluster the two cases. More over, GMM
is computationally extensive when compared
to k-meons.

Alglomerative Clustering is bottom-top approach.

Hence from # datapoints to 2 clusters it will take large no of iderations. Hence, k-means a is better than aglomerative clustering.

Since, that do not have outliers in the final set (as outliers are semoved by ICBR thoushold). Hence, DBSCAN will not outperform K-means dustering in this case.

2);i) For evaluating the clustering algorithm
we can use the which of the algorithm (1c-mass,

CIMM, hierarchical, DBSCAN) was made a tighter

clusterie cost function is minimum.

1. \(\sum_{i=1}^{1} \) \(\sum_{i=1}^{

Note: Multiple solutions are also allowed for this question.

2)(iii)
Multiple solutions are allowed

(iv) Multiple solutions are allowed.

MARKING SCHEME

1) a) (i) 1 Marks

(ii) 1 Marls

(iii) 1 Marles for firding minimum using code 1 Merls for pros e cons

(iv) 1 Marla

1) 6) (1 Marle) for difference between regression e k-means

(1 Maroles) for comment on cost for

(2 Mars) for illustrating using code.

(1 Marts) Propose method & justification.

2) (i) [Model] t-SNE + prolipsocessing (0.5) (0.5)

I Mayor for Justification.

(ii) (1 Marks) for evaluation metric
1 Marks for coding

(iii) Poetsocessing - (Marla)

To show degradation - (IMordes)

Final proposal - (1 Marles)

(ji) Block diagram & equation (2 mara) (ode validation (1 mares)