#### CS6011: Kernel Methods for Pattern Analysis

Even 2015

## Assignment II Report

15/03/2015

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# 1 Objective

The objective of this assignment is to understand the application of MLFFNN in both classification and regression tasks.

- 1. Under classification tasks, the performance of MLFFNN is studied in following 4 cases
  - (a) **Linearly separable classes** Here, the performance of MLFFNN with 1 or 2 hidden layers is compared with Bayes classifier using GMM model and perceptron. Number of classes is 3.
  - (b) **Non-linearly separable classes** Here, the performance of MLFFNN with 1 or 2 hidden layers is compared with Bayes classifier using GMM model. Number of classes is 2.
  - (c) **Overlapping classes** Here, the performance of MLFFNN with 1 or 2 hidden layers is compared with Bayes classifier using GMM model. Number of classes is 3.
  - (d) **Image data** Here, the performance of MLFFNN with 2 hidden layers is compared with Bayes classifier using GMM model and perceptron. Number of classes is 5.
- 2. Under regression tasks, function approximation of Univariate and Bivariate dataset are experimented with MLFFNN and RBF models.
  - (a) To find the best performing MLFFNN model with 1 hidden layers for the Univariate dataset
  - (b) To find the best performing MLFFNN model with 1 or 2 hidden layers for the Bivariate dataset
  - (c) To find the best performing RBF model for both Univariate and Bivariate datasets and to analyse the effect of regularisation parameter and model complexity on the performance of RBF model.

#### 2 Classification tasks

## 2.1 Linearly separable classes

The given data includes 3 linearly separable classes. Visualisation of the training data is as in Figure  $\,1\,$ 

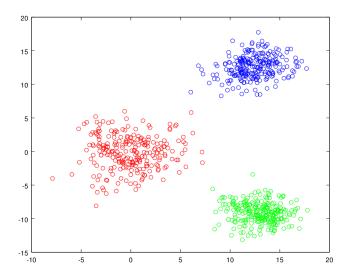


Figure 1: Linearly separable classes

#### 2.1.1 Bayes classifier using Gaussian model

Bayes classifier involves Bayesian reasoning which provides a probabilistic approach to inference. It is based on certain assumptions like the underlying data distribution follows a particular probability distribution and enables to make optimal decisions in classification tasks by making use of the assumed probabilities together with observed data. As per the bayes rule,

$$P(\omega_j|x) = \frac{p(x|\omega_j)P(\omega_j)}{p(x)}$$

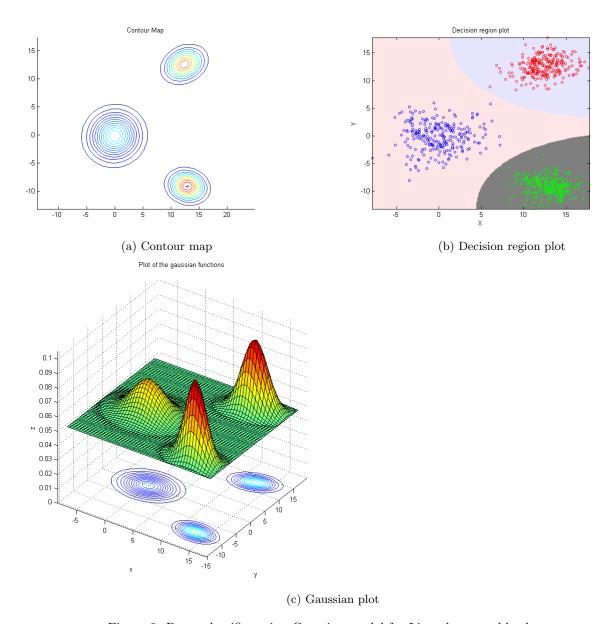
The above formula calculates the posterior probability that a particular point  $\bar{x}$  belongs to class  $\omega_j$ . We will assign  $\bar{x}$  to the class j whose posterior probability is the maximum.

Coming on to the right hand side of this equation,  $P(\omega_j)$  is same for all classes in our data set because each class has same number of training examples and we do not know anything about the real world prior distribution. Also, denominator is ignored as it is same for all classes for a particular value of x, our decision is solely dependent on  $p(x|\omega_j)$ .

For this assignment, we have considered gaussian distribution function for deciding  $p(x|\omega_j)$ . This function is dependent on mean and covariance of the data and is decided based on that. This function is given by

$$\frac{1}{\sigma\sqrt{2\pi}}e^{-(x-\mu)^2/2\sigma^2}$$

In order to deal with calculations on very small values of probabilities, we take the log of likelihood function and then calculate the probabilities.



 $\label{thm:classifier} \mbox{ Figure 2: Bayes classifier using Gaussian model for Linearly separable classes}$ 

The confusion matrix for both validation and test data are shown below:

	class1	class2	class3
class1	150	0	0
class2	0	150	0
class3	0	0	150

Table 1: Validation data

	class1	class2	class3
class1	100	0	0
class2	0	100	0
class3	0	0	100

Table 2: Test data

- One of the three classes has a circular contour which suggests independence of features in that class i.e. diagonal covariance matrix with all variances same
- 3 gaussians are enough to classify the data (one for each class)
- the means of the three classes are well separated from each other
- the covariance matrices of the three classes are different implies a non-linear decision boundary

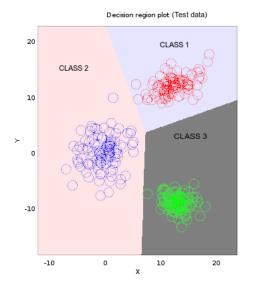
#### 2.1.2 Perceptron

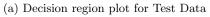
The perceptron is a type of artificial neural networks. It can be seen as the simplest kind of feedforward neural network: a linear classifier. We experimented with both pattern mode of learning and batch mode of learning.

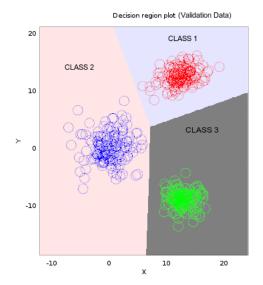
#### Pattern mode

- Initialize the weight vector  $w < w_0, w_1, \dots > to 0$ .
- Repeat
- Alter the weights for every misclassified point in training data.  $w(m+1) = w(m) + \eta^* y^* x(m)$
- till all the points are properly classified.

We trained the perceptron on the training data provided to us with  $\eta = 0.1$ . The decision region plots for the validation and test data is given below.







(b) Decision region plot for Validation data

Figure 3: Classification using Perceptron

The confusion matrix for the test data and validation data is given below:

	Α	В	С
Α	150	0	0
В	0	150	0
С	0	0	150

(a) Validation Data

	Α	В	С
Α	100	0	0
В	0	100	0
С	0	0	100

(b) Test data

Figure 4: Confusion matrix

The accuracy of the learned model is 100%.

#### Batch Mode Learning:

In this variant of perceptron learning we scan all of the training examples once and then update the weight at the end one epoch. Cumulative changes of all misclassified examples is done at the end.

#### Algorithm:

- Initialize the wight vector w to 0.
- Repeat for each epoch
- $\triangle w(1) = 0$
- Repeat for each training data sample
- If the sample is misclassified
- $\triangle w(l) = \triangle w(l) + x*y$

- $\bullet$  Endif
- $w(l+1) = w(l) + \eta * \triangle w(l)$
- End epoch

The perceptron weights were learned with  $\eta=0.1$ . The decision plots for the test data and the validation data after learning perceptron in batch mode are:

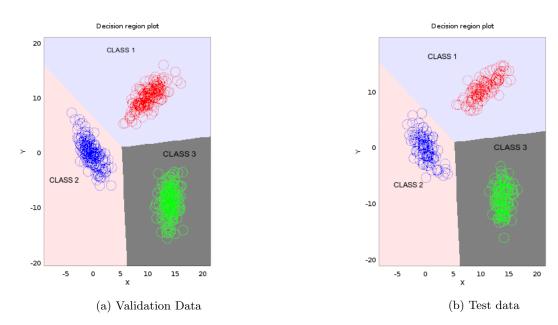


Figure 5: Decision region plots

Confusion matrix for the test and train data is as follows:

	Α	В	С
Α	100	0	(
В	0	100	C
С	0	C	100
		(a) Te	st data

Figure 6: Confusion matrix

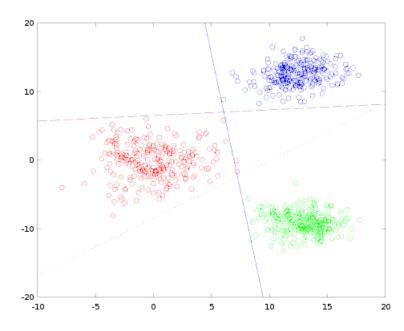
The accuracy of the learned model is 100%.

Learning Rate	Epocs	Epocs	Epocs	Accuracy
	Class A	Class B	Class C	Testdata
0.1	500	8	5	100
0.2	500	3	6	99.67
0.3	500	18	8	100
0.4	500	7	16	100
0.5	500	15	19	100
0.6	500	16	17	100
0.7	500	9	9	100
0.8	500	8	8	100
0.9	500	9	8	99.67

Table 3: Learning rate analysis with respect to number of epochs for pattern mode

Learning Rate	Epocs	Epocs	Epocs	Accuracy
	Class A	Class B	Class C	Testdata
0.1	500	34	31	99.67
0.2	500	34	31	100
0.3	500	8	5	100
0.4	500	8	5	99.67
0.5	500	34	31	99.67
0.6	500	34	31	99.67
0.7	500	34	31	99.67
0.8	500	34	31	99.67
0.9	500	34	31	100

Table 4: Learning rate analysis with respect to number of epochs for batch mode



(a) Separating line obtained by the perceptron learn on training data

- Batch mode takes more number of epochs to converge
- In pattern mode, we can observe that the perceptron converges faster with increase in learning rate
- The perceptron that is learned for the class 1 Vs rest is not converging even at 500 epochs. This is because there is one data point in training set of Class 1 that lies on the boundary. Every time we alter the weights near the separating line, at least one sample is misclassified.

#### 2.1.3 MLFFNN

Multilayer feedforward neural network consists of multiple hidden layers which are connected in feedforward way. There are three main layers - input layers, hidden layers and output layers. The number of hidden layers decides the complexity of the decision surface.

We used gradient descent method to train linearly separable data. The data is trained with one hidden layer. The number of hidden nodes are chosen by cross validation. The plot for the mean square error versus number of hidden nodes is shown in Figure 7.

#### Inferences

- Minimum MSE is obtained for 3 hidden nodes
- Fixed the number of hidden nodes as 3

In order to choose the learning rate, we plotted the MSE versus number of epocs for different learning rates. The Figure 8 depicts the plots for learning rate 0.1 and 0.9. For these plots we used hyperbolic tangent activation function. The plot for logistic sigmoid activation function is shown in Figure 9.

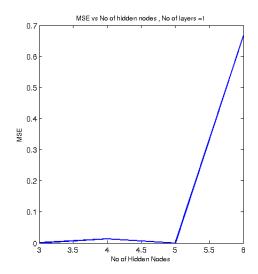
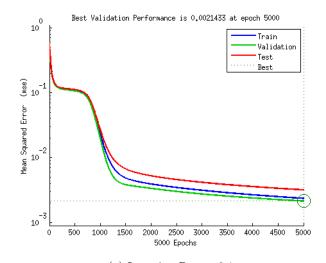
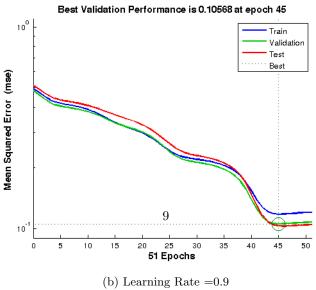


Figure 7: Linearly separable classes







( )

Figure 8: Learning Rate analysis for hyperbolic tangent activation function

- No oscillations observed for learning rate=0.1
- $\bullet$  At learning rate =0.9, we can see oscillations
- Minimum MSE achieved at learning rate =0.1

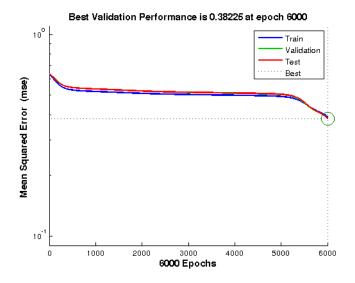


Figure 9: At Learning Rate =0.1 for Logistic activation function

#### Inferences

- MSE is more compared to hyperbolic tangent activation results
- Hyperbolic tangent function is faster than sigmoid function

Finally we chose the model with one hidden layer and 3 hidden nodes, hyperbolic tangent activation function and learning rate as 0.1. The network view is shown in Figure 10.

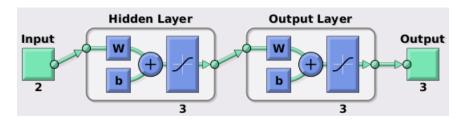


Figure 10: Neural Network View

The decision plot obtained for the given data using the mentioned model parameters is shown in Figure 11.

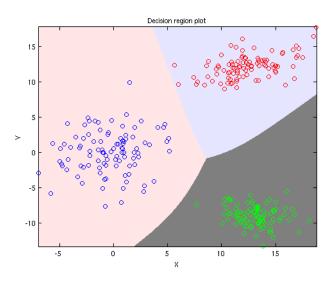


Figure 11: Decision region plot

The model gave 100% accuracy. The corresponding confusion matrix for validation data and test data are shown in Figure 12.



Figure 12: Confusion Matrix

• 100% accuracy obtained for both validation data and test data.

#### 2.1.4 Comparison of performance of different models

## Accuracy Comparison across different models

Method	Validation	Test
	Accuracy (%)	Accuracy (%)
Bayes Classifier	100	100
Perceptron	100	100
MLFFN	100	100
1 hidden layer		

#### Inferences

• All models are working well for linearly separable data

## 2.2 Non-linearly separable classes

The given data includes 2 linearly separable classes. Visualisation of the training data is as in Figure 13

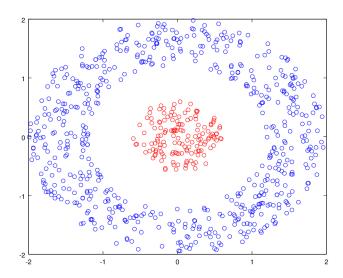
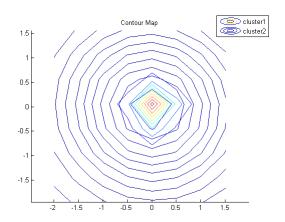
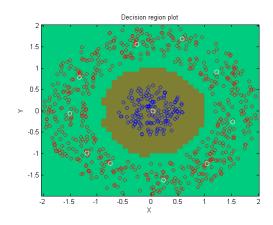


Figure 13: Non-linearly separable classes

#### 2.2.1 Bayes classifier using GMM and Gaussian





- (a) Contour map of GMM with 1 gaussian for each class
- (b) Decision region plot of GMM with 10 gaussians for class 1 and 1 for class 2

Figure 14: Bayes classifier using GMM for Non-linearly separable classes

The confusion matrices for validation and testing data are as shown below:

	class1	class2
class1	90	0
class2	0	360

Table 5: Validation data

	class1	class2
class1	60	0
class2	0	240

Table 6: Test data

#### Inferences

- GMM using 10 clusters (as highlighted in the Figure 14) gives 100 percent accurracy.
- A single gaussian also gives 100% accuracy in this case. This is because of the contour of gaussian 1 is concentric to that of gaussian 2. However, this might not give good accuracy on all non linearly separable datasets.

#### 2.2.2 MLFFNN

The data is trained using gradient descent method with one hidden layer. The number of hidden nodes are decided using cross validation The Figure 18 shows the plot of mean square error versus number of hidden nodes.

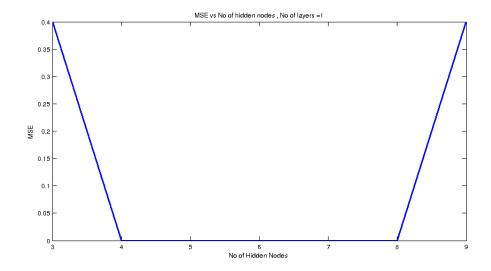


Figure 15: MSE vs No of Hidden nodes

• Minimum mean square error obtained for 4 hidden layers

Then the number of epochs versus MSE is plotted for learning rate 0.1 and 0.9. The plots are shown in Figures 16 and 17.

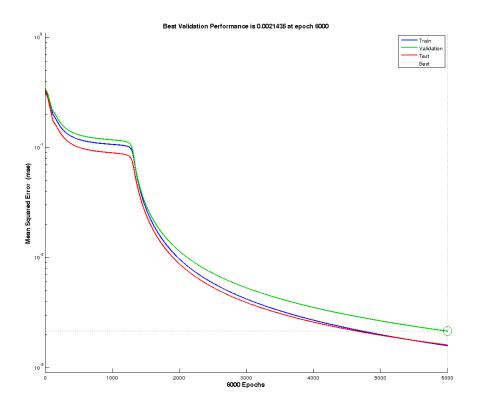


Figure 16: No of epochs vs MSE at learning rate =0.1

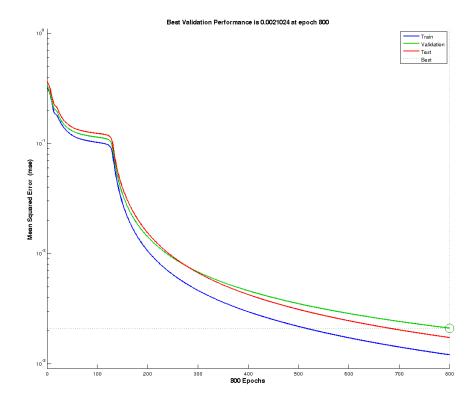


Figure 17: No of epochs vs MSE at learning rate =0.9

- No oscillations are found in both the cases.
- $\bullet\,$  Fixed learning rate as  $0.1\,$

We used hyperbolic tangent function as activation function in both hidden layer and output layer. The view of the network is shown in Figure 19.

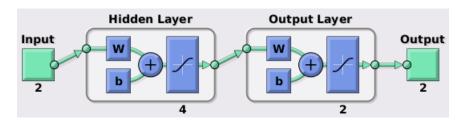


Figure 18: View of Neural network

The decision plot obtained for the given data using the mentioned model parameters is shown in Figure 19.

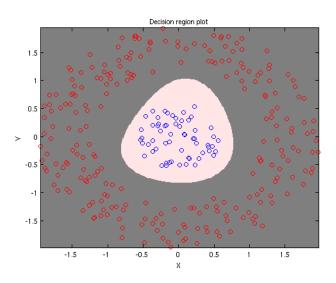


Figure 19: Decision region plot

The model gave 100% accuracy. The corresponding confusion matrix for validation data and test data are shown in Figure 20.



Figure 20: Confusion Matrix

 $\bullet$  100% accuracy obtained for both validation data and test data.

# 2.2.3 Plots of Outputs of each of the hidden nodes and output nodes Using 1 hidden layer for no of epochs = 1

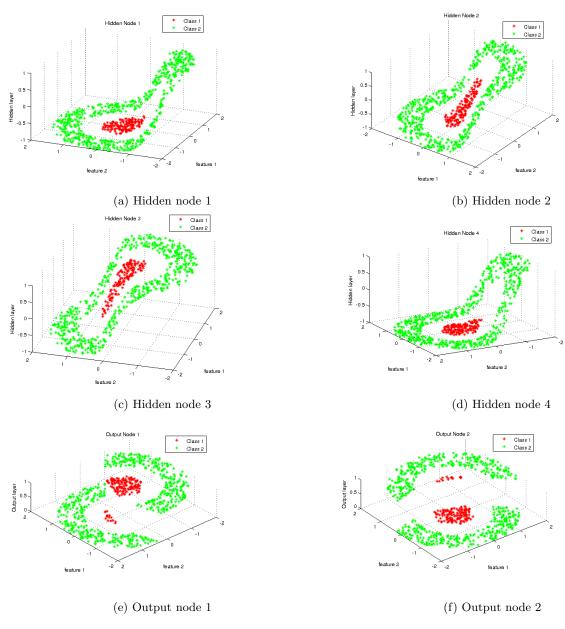


Figure 21: No of Epochs =1

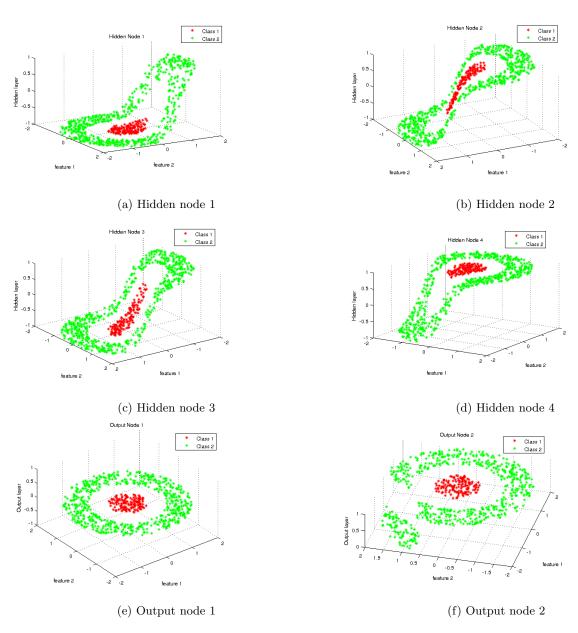


Figure 22: No of Epochs =2

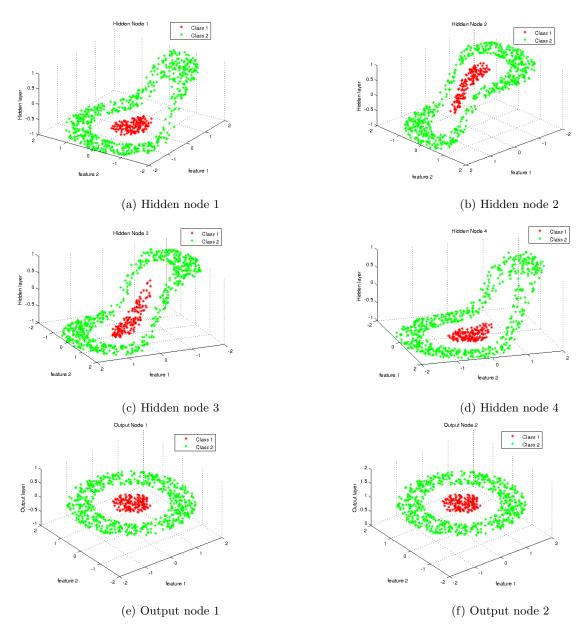


Figure 23: No of Epochs =10

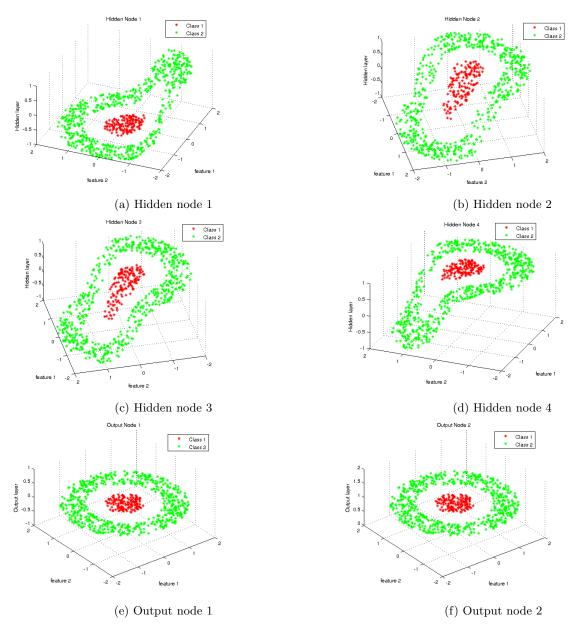


Figure 24: No of Epochs =50

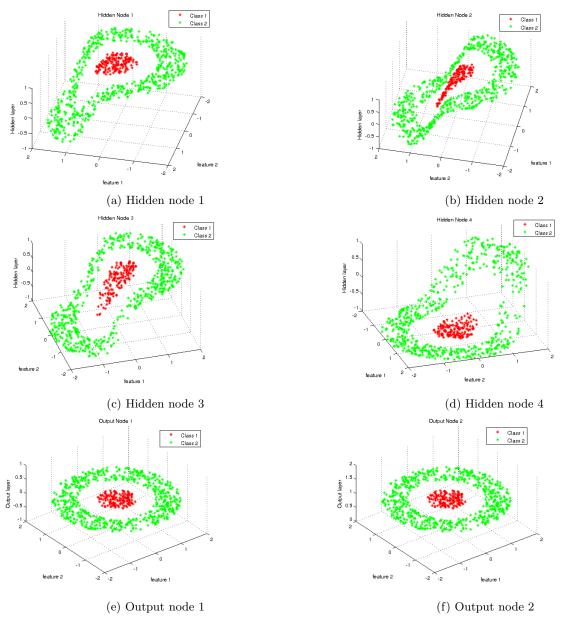


Figure 25: No of Epochs =100

#### Inferences

- Non linear tangent activation function results an S-shaped surface for each hidden node plots.
- From the output nodes of each epochs we can observe that as the number of epochs increases, the classification

accuracy is increasing

- At 100 epochs all data points are classified into one class(majority). This is due to the class imbalance in the given data.
- With more number of epochs (500), model classifies correctly

## Using 2 hidden layer for no of epochs = 1

Our best model uses 4 hidden nodes in the first hidden layer and 3 hidden nodes in the second hidden layer.

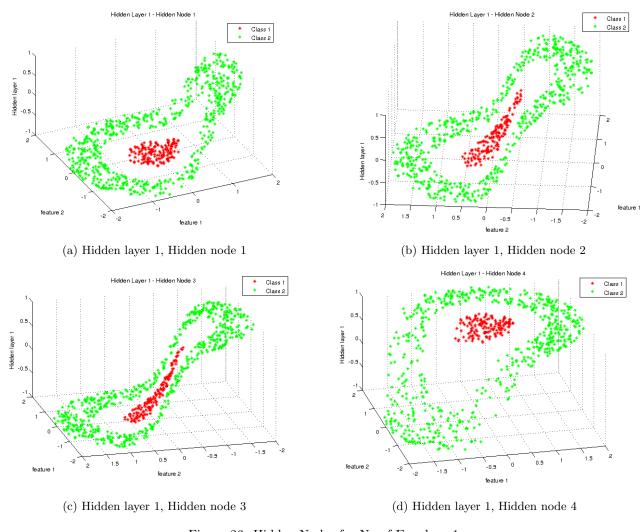
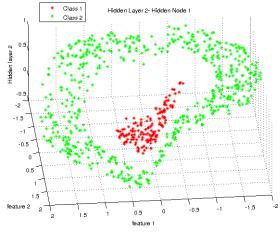
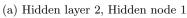
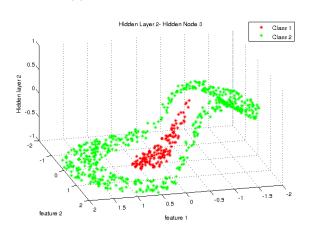


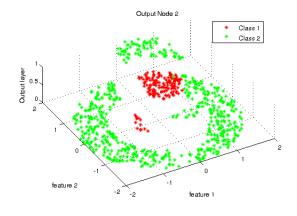
Figure 26: Hidden Nodes for No of Epochs =1



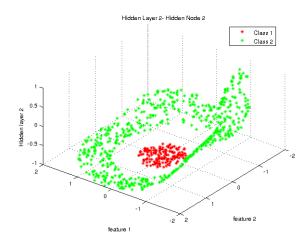




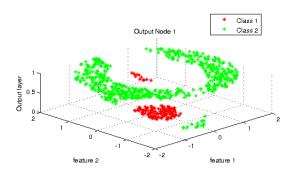
 $\left( c\right)$  Hidden layer 2, Hidden node 3



(e) Output node 2



(b) Hidden layer 2, Hidden node 2



(d) Output node 1

Figure 27: Hidden and Output nodes for No of Epochs =1

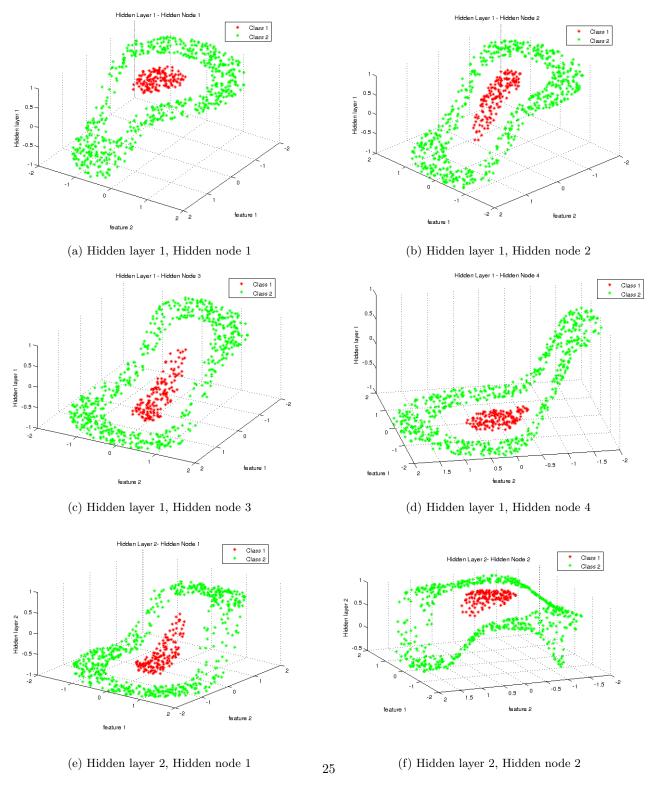
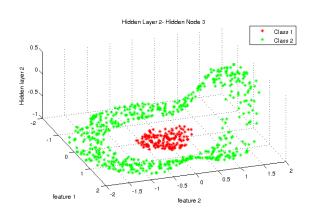
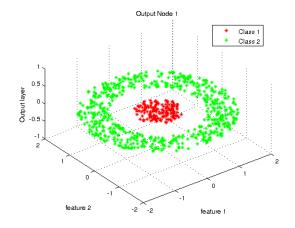
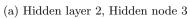
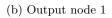


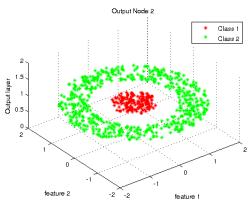
Figure 28: Hidden Nodes for No of Epochs =2











(c) Output node 2

Figure 29: Output nodes for No of Epochs =2

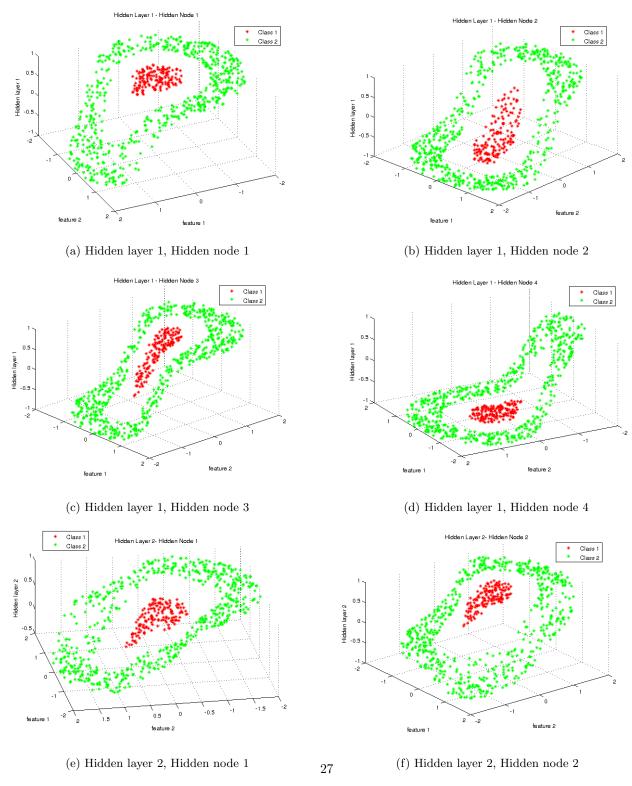
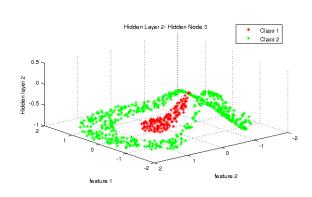
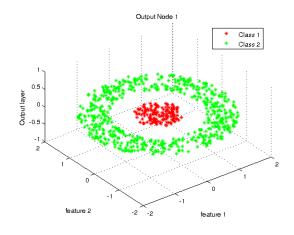
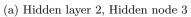
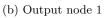


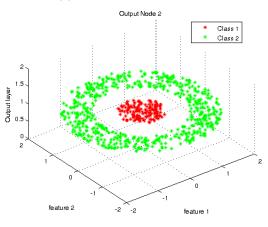
Figure 30: Hidden Nodes for No of Epochs =10  $\,$ 











(c) Output node 2

Figure 31: Output nodes for No of Epochs =10

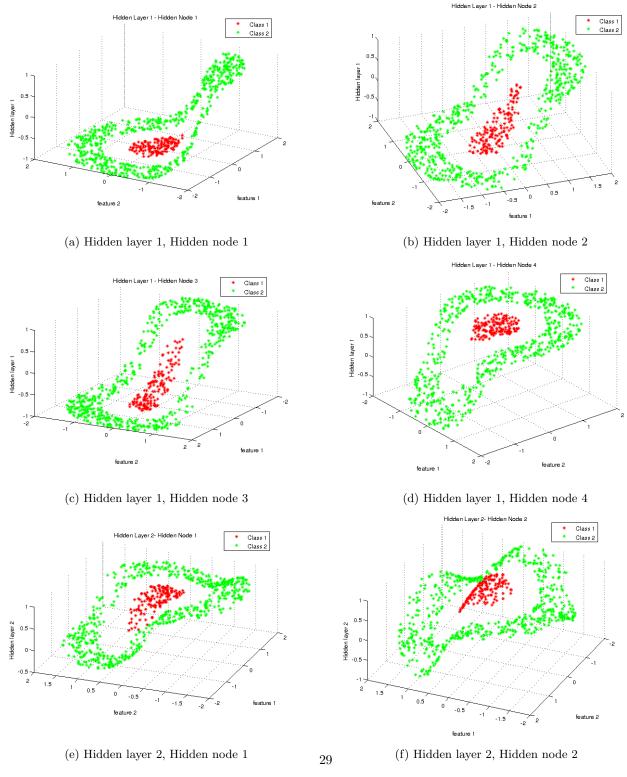


Figure 32: Hidden Nodes for No of Epochs =50

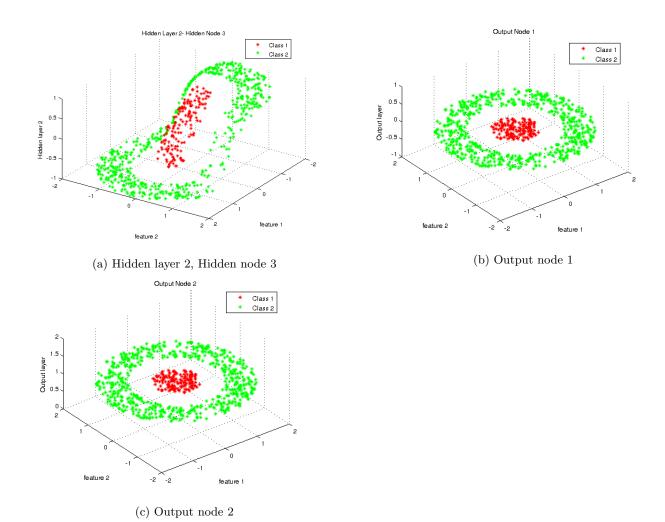


Figure 33: Output nodes for No of Epochs =50

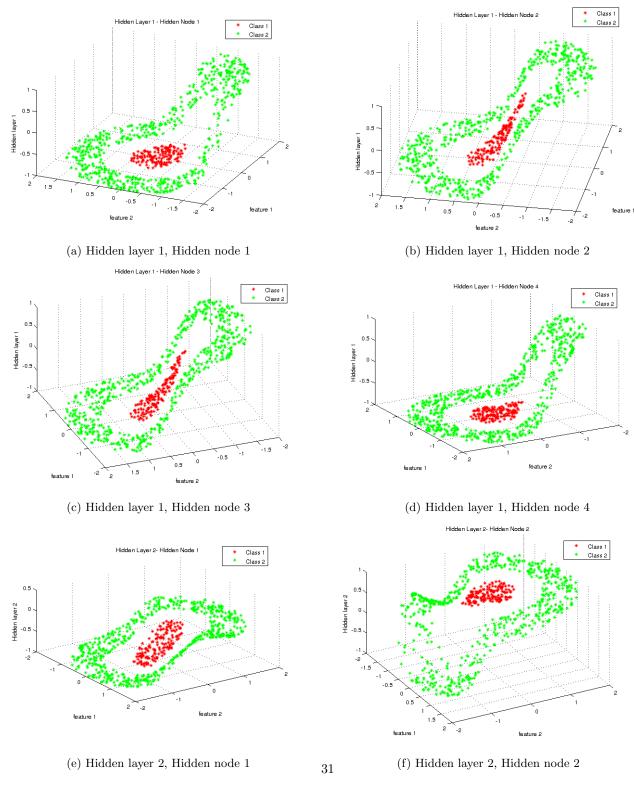
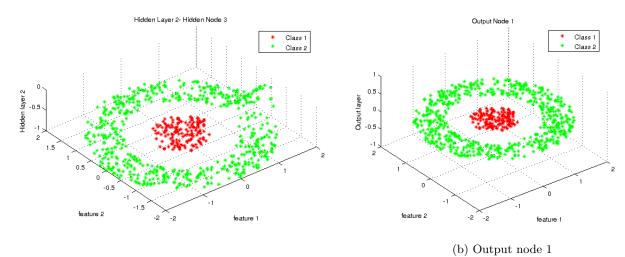
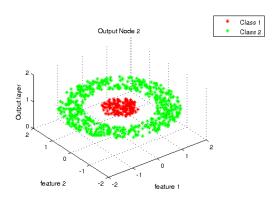


Figure 34: Hidden Nodes for No of Epochs =100



(a) Hidden layer 2, Hidden node 3



(c) Output node 2

Figure 35: Output nodes for No of Epochs =100

### Inferences

• For number of epochs=2, all data points get classified into one class

## 2.2.4 Comparison of performance of different models

Accuracy Comparison across different models

Method	Validation	Test	
	Accuracy (%)	Accuracy (%)	
Bayes Classifier	100	100	
MLFFN	100	100	
1 hidden layer			

# 2.3 Overlapping classes

The given data includes 3 linearly separable classes. Visualisation of the training data is as in Figure 36

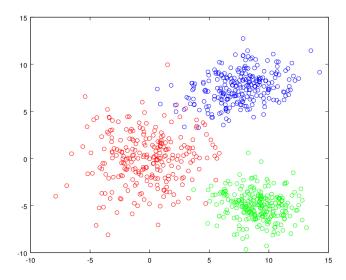


Figure 36: Overlapping classes

## ${\bf 2.3.1}\quad {\bf Bayes\ classifier\ using\ Gaussian\ model}$

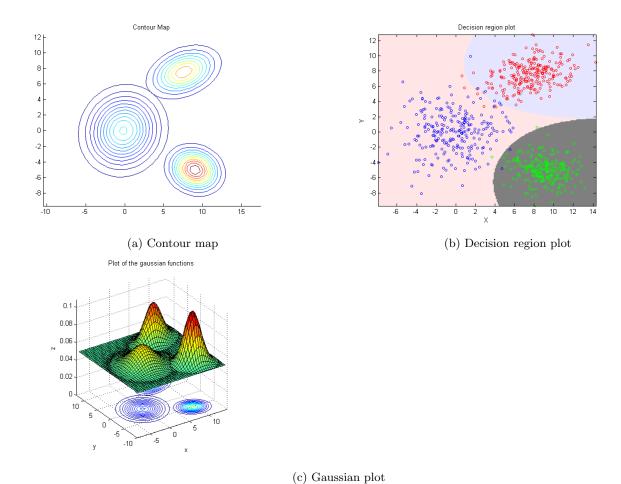


Figure 37: Bayes classifier using Gaussian model for Overlapping classes

The confusion matrices for validation and testing data are as follows:

	class1	class2	class3
class1	146	3	1
class2	1	149	0
class3	1	0	149

Table 7: Validation data

	class1	class2	class3
class1	97	1	2
class2	2	98	0
class3	0	0	100

Table 8: Test data

- 3 gaussians are used to model the given classes
- Accuracy obtained on validation data is 98.67%
- Accuracy obtained on test data is 98.33%

#### 2.3.2 MLFFNN

For overlapped data, we used two hidden layers. The number of hidden nodes in each hidden layer are found using cross validation. The Figure 38 is plotted to decide the number of hidden nodes in first hidden layer. For second hidden layer we obtained the number of hidden nodes as 3.

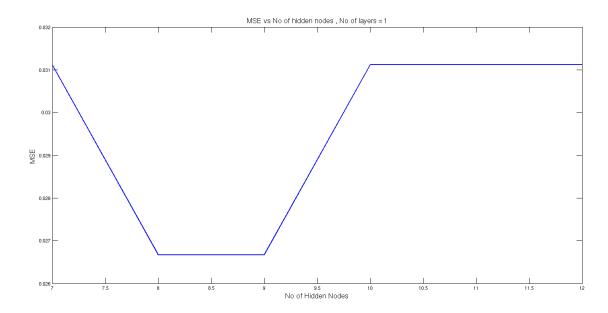


Figure 38: MSE vs Number of hidden nodes

#### Inferences

• Minimum MSE obtained for 8 hidden nodes in hidden layer 1 and 3 hidden nodes in hidden layer 2

We fixed the learning rate as 0.4 for this experiment (using cross validation). Also, we used generalized delta rule for weight update by using *traingdm*. The momentum factor we used is 0.8. The network view of our model is shown in Figure 39.

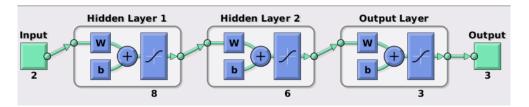


Figure 39: View of network

The decision plot obtained for the given data using the mentioned model parameters is shown in Figure 40.

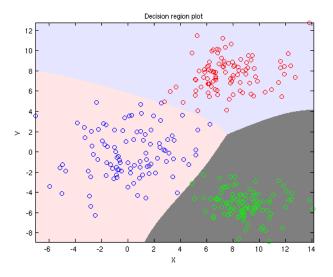


Figure 40: Decision region plot

The model gave 98% accuracy. The corresponding confusion matrix for validation data and test data are shown in Figure 41.

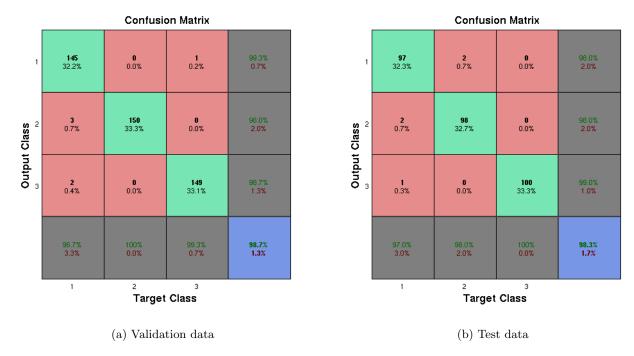


Figure 41: Confusion Matrix

- $\bullet~98.7\%$  accuracy obtained for validation data
- 98.3% accuracy obtained for test data

## 2.3.3 Comparison of performance of different models

### Accuracy Comparison across different models

Method	Validation	Test	
	Accuracy (%)	Accuracy (%)	
Bayes Classifier	98.6	98.3	
MLFFN	98.7	98.3	
1 hidden layer			

# Inferences

 $\bullet\,$  Both are performing similar

# 2.4 Image data

# 2.4.1 Bayes classifier using GMM

We used GMM on the image dataset with the following number of clusters.

class 1 : 15 class 2 : 15 class 3 : 25 class 4 : 25 class 5 : 25

We obtained 62% accuracy on the validation and test data. Confusion matrices are as shown below:

	1	2	3	4	5
1	20	0	0	0	0
2	6	18	1	0	3
3	6	0	8	0	1
4	8	5	0	2	1
5	4	0	1	1	13

Table 9: Validation Data

	1	2	3	4	5
1	32	7	0	0	1
2	16	35	1	1	4
3	1	0	25	0	3
4	12	10	0	4	5
5	11	2	0	0	25

Table 10: Test data

### Inferences

The classes 2,4,5 are frequently confused with those in classes 1.

## 2.4.2 MLFFNN with 2 hidden layers

We used Levenberg-Marquardt optimization method for training which was giving more performance than gradient descent method. The activation function used is hyperbolic tangent function. The learning rate is chosen as 0.4 using cross validation. Two hidden layers are used. The plot for number of hidden nodes in hidden layer 1 versus MSE is shown in Figure 43.

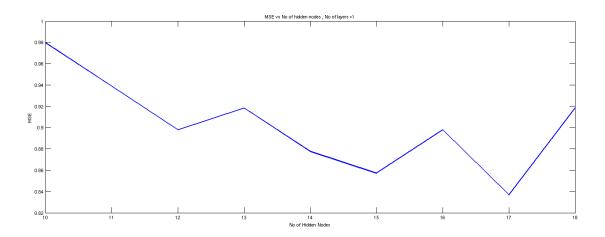


Figure 42: MSE vs No of hidden nodes in hidden layer  $1\,$ 

• The minimum error obtained for 17 hidden nodes.

By cross validation we chose the number of hidden nodes in the hidden layer 2 as 3.

Then we trained the model using these parameters and the confusion matrix for validation data and test data are shown in Figure 43.

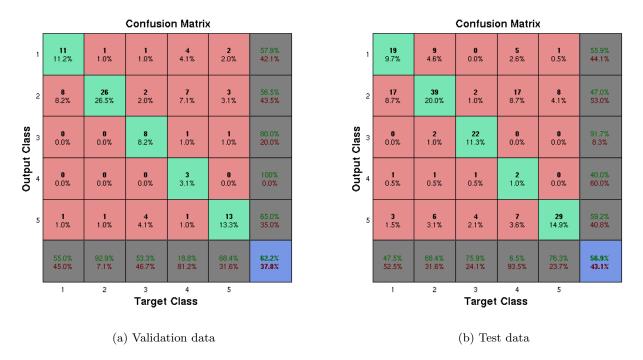


Figure 43: Confusion Matrix

- Obtained 62.2% accuracy for validation data and 56.9% accuracy for test data
- The performance is poor compared to other datasets, since this is a real time data.

### 2.4.3 Comparison of performance of different models

#### Accuracy Comparison across different models

Method	Validation	Test	
	Accuracy (%)	Accuracy (%)	
Bayes Classifier	62	62	
MLFFN	62.2	56.9	
2 hidden layer			

#### Inferences

- Bayesian is performing better in test data
- We are not getting better accuracy because this is real time data

# 3 Regression tasks

# 3.1 MLFFNN on Univariate Dataset

The underlying function of univariate dataset is  $\cos^2(2\pi x)$ . A gaussian noise with zero mean is added. For this experiment we generated 5000 samples for training data, 1000 samples for validation data and 500 samples for test data. We plotted the values of mean squared error(MSE) on training data, validation data and test data across varying number of hidden nodes. The plot is shown in Figure 44.

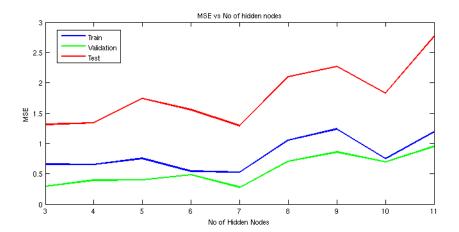


Figure 44: MSE versus No of hidden nodes

 $\bullet$  The minimum validation error obtained for No of hidden nodes=12

The plot across model output and target output for three sets data are shown in Figure 45.

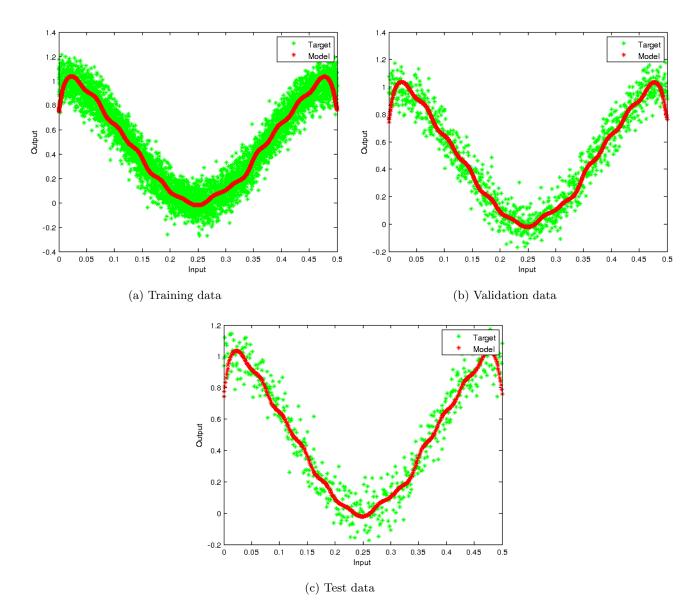


Figure 45: Model output and target output for training data, validation data and test dat

## Inferences

 $\bullet$  The approximated function is somewhat fitting the underlying function

The scatter plots with target output on x axis and model output on y axis for training data, validation data and test data are shown in Figure 46.

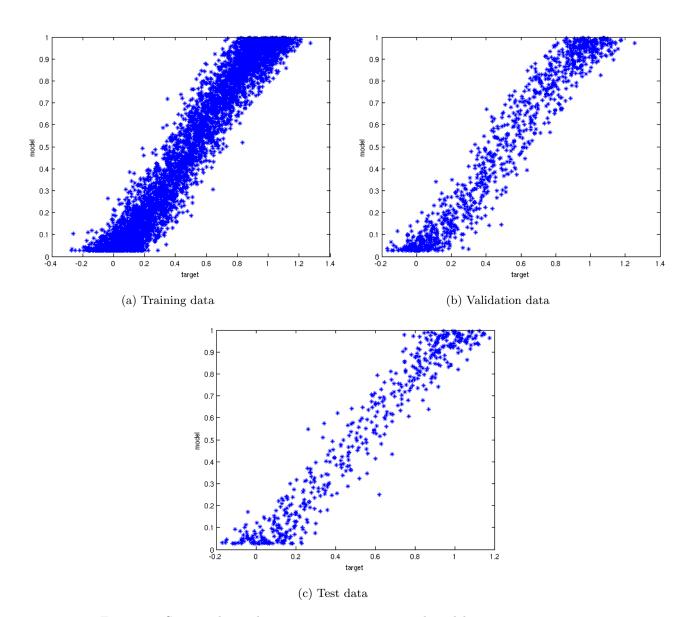


Figure 46: Scatter plot with target output as x axis and model output as y axis

#### Result

Error	MSE	
Training	0.0086	
Validation	0.0081	
Testing	0.0085	

#### Inferences

- The scatter plots for all the three datasets are close to y=x line.
- The error is similar to regression tasks.

### 3.2 MLFFNN on Bivariate Dataset

We experimented the bivariate datasets for various training data size. The plots obtained for training data size 100 and 2000 are shown in this report. We trained the data using gradient descent method and Levenberg-Marquardt optimization method. The number of hidden nodes we chose using cross validation and we experimented for one hidden layer and two hidden layer. We used hyperbolic tangent function as activation function in hidden layers and linear activation function in output layer. The learning rate fixed through cross validation and it is fixed as 0.1.

The plot of the values of MSE on training data, validation data and test data for different hidden nodes are shown in Figure 47.

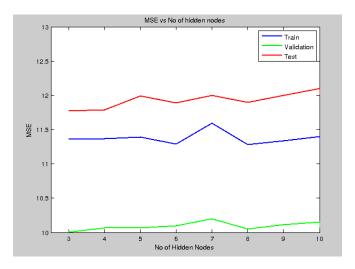


Figure 47: MSE vs No of hidden nodes

#### Inferences

• The best performance is at no of hidden nodes =3

So we chose number of hidden nodes in first hidden layer as 3 and second hidden layer as 2.

The plot across model output and target output for three sets data with training data size as 2000 and one hidden layer are shown in Figure 48.

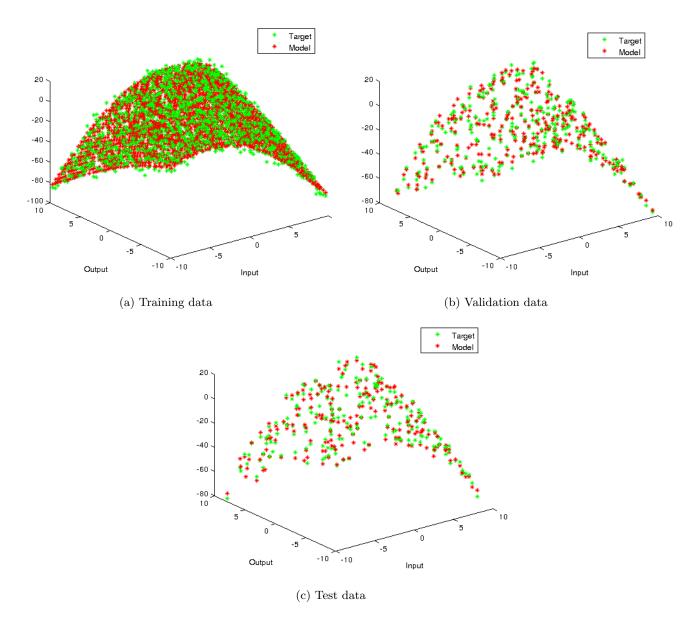


Figure 48: Model output and target output for training data, validation data and test data with training data size=2000

The plot across model output and target output for three sets data with training data size as 100 and one hidden layer are shown in Figure 49.

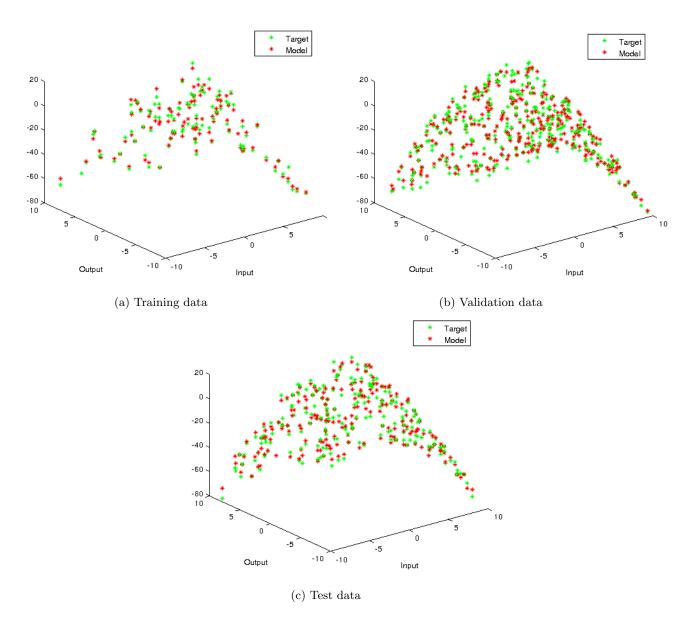


Figure 49: Model output and target output for training data, validation data and test data with training data size=100

The plot across model output and target output for three sets data with training data size as 2000 and two hidden layer are shown in Figure 50.

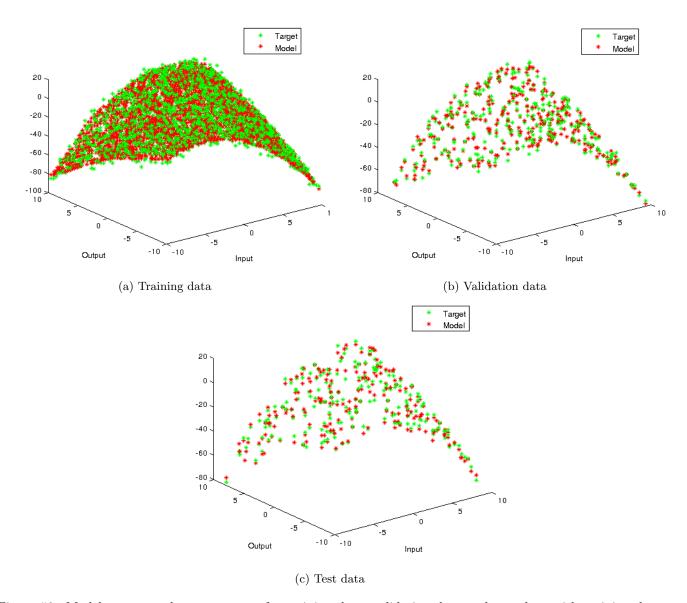


Figure 50: Model output and target output for training data, validation data and test data with training data size=2000 obtained using two hidden layer

The Figures 48, 49 and 50 are obtained using Levenberg-Marquardt optimization method. The plots for the same using gradient descent method for training data size 2000 is shown in Figure 51.

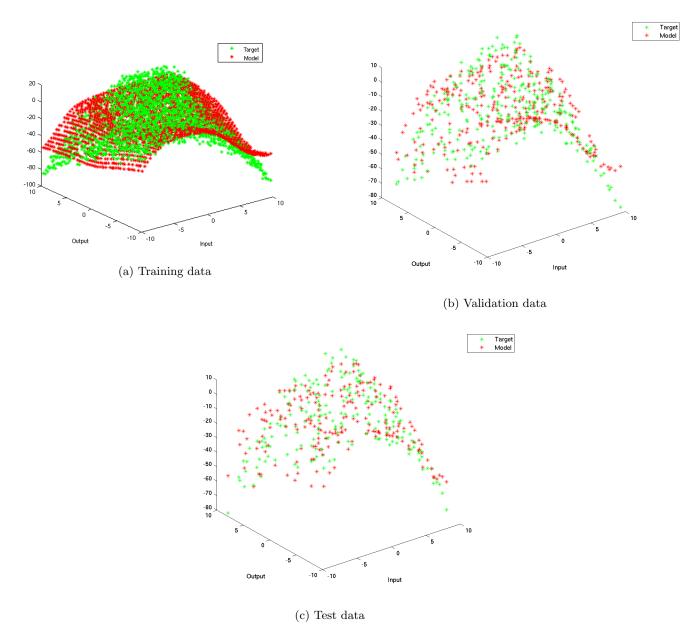


Figure 51: Model output and target output for training data, validation data and test data with training data size=2000 obtained using gradient descent method

• Levenberg-Marquardt optimization method performs well than gradient descent method

The scatter plot with target output on x-axis and model output on y-axis for three sets data with training data size as 2000 and one hidden layer are shown in Figure 52.

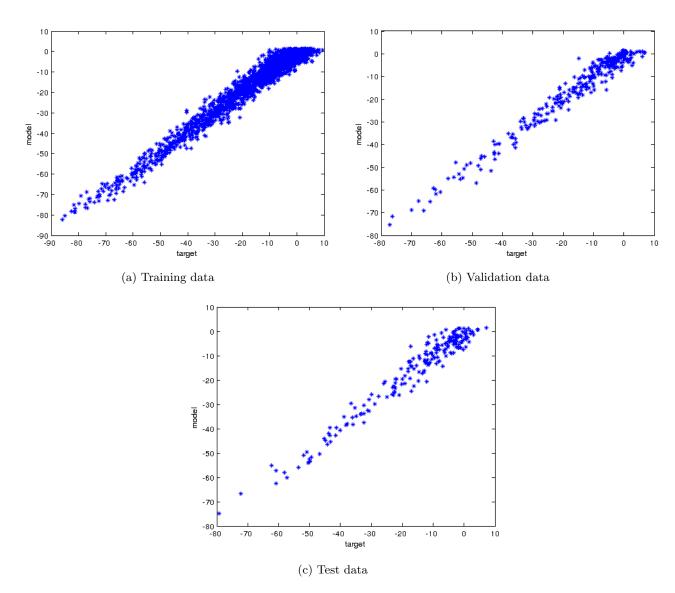


Figure 52: Scatter plot with target output on x-axis and model output on y-axis with training data size=2000

The scatter plot with target output on x-axis and model output on y-axis for three sets data with training data size as 100 and one hidden layer are shown in Figure 53.

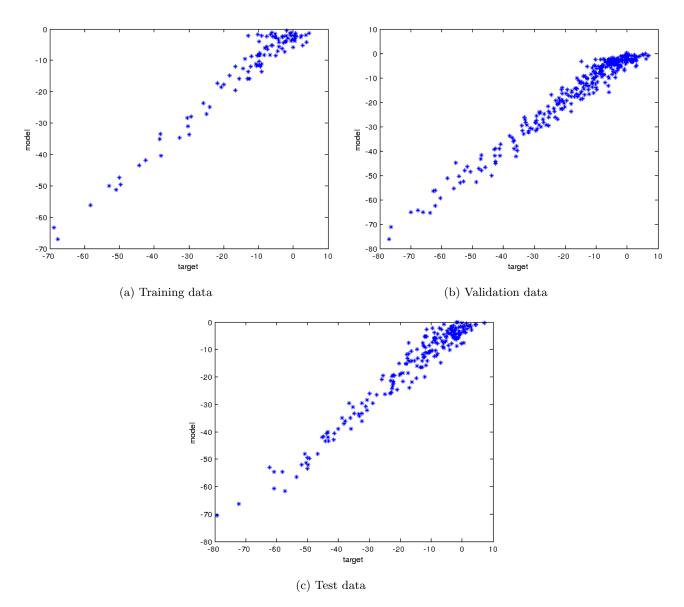


Figure 53: scatter plot with target output on x-axis and model output on y-axis for training data, validation data and test data with training data size=100

The scatter plot with target output on x-axis and model output on y-axis for three sets data with training data size as 2000 and two hidden layer are shown in Figure 54.

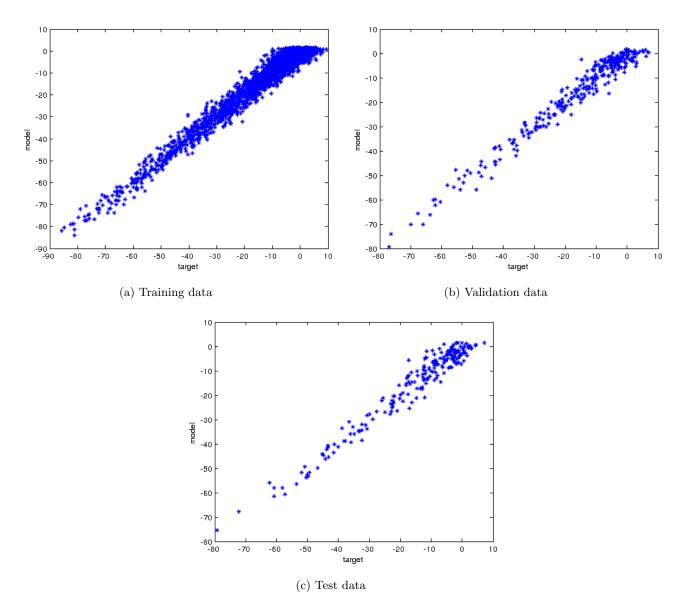


Figure 54: scatter plot with target output on x-axis and model output on y-axis for training data, validation data and test data with training data size=2000 obtained using two hidden layer

The Figures 52, 53 and 54 are obtained using Levenberg-Marquardt optimization method. The plots for the same using gradient descent method for training data size 2000 is shown in Figure 55.

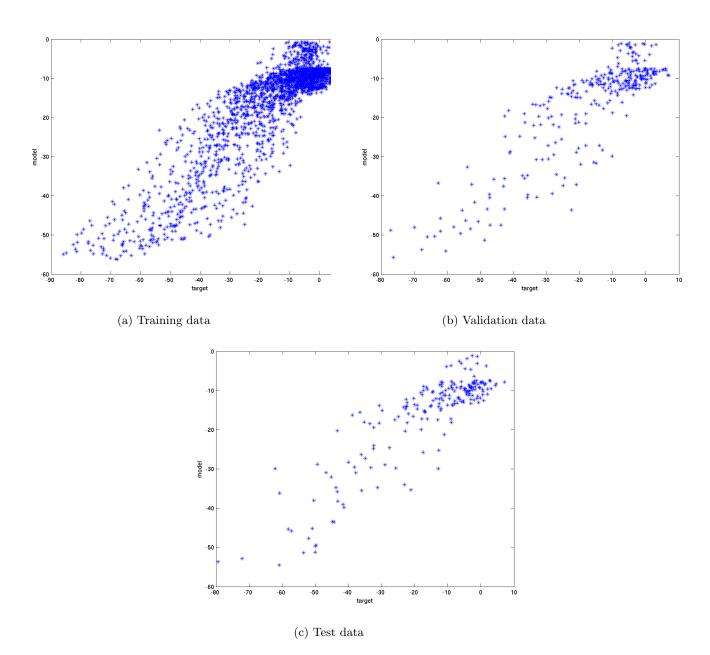


Figure 55: scatter plot with target output on x-axis and model output on y-axis for training data, validation data and test data with training data size=2000 obtained using gradient descent method

#### Result

Error	MSE for N=2000	MSE for N=2000	MSE for $N=2000$	MSE for N=100
	One hidden layer	Two hidden layer	Using Gradient descent	One hidden layer
Training	11.46	11.22	81.13	11.33
Validation	10.06	10.06	78.60	11.74
Testing	11.82	12.05	75.47	12.69

#### Inferences

- The points in the scatter plots are across 45 degree inclination of the axes indicating that the modeled data follow the target data very well
- The plots obtained for gradient decent method based training doesn't show much inclination as other method do
- Error is less compared to regression models
- MSE obtained for the model using gradient descent method is too high compared to Levenberg-Marquardt method

### 3.2.1 Plots of Outputs of each of the hidden nodes and output nodes

The best model obtained for one hidden layer contains 3 hidden nodes. The plot of outputs of each of the hidden nodes and output nodes for various number of epochs are shown below. These plot are generated for training data set size 2000.

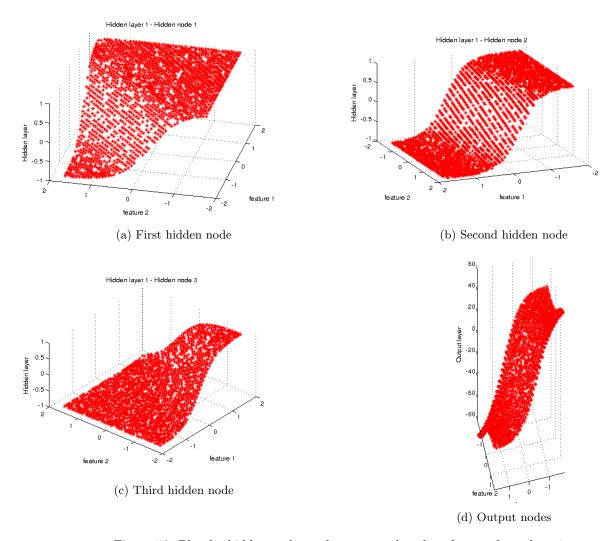


Figure 56: Plot for hidden nodes and output node values for no of epochs =1

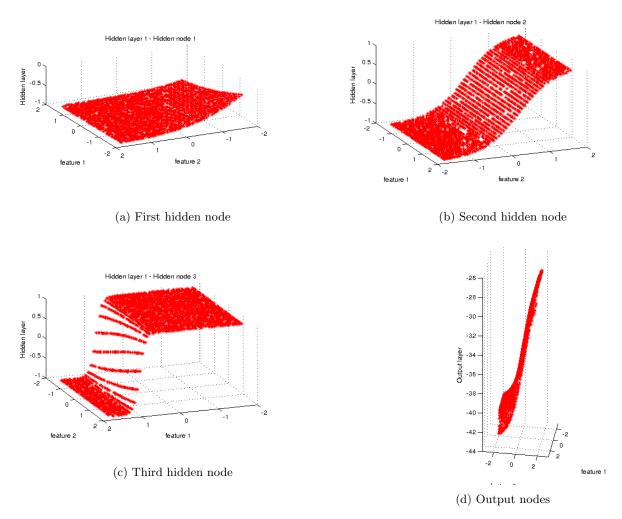


Figure 57: Plot for hidden nodes and output node values for no of epochs =2

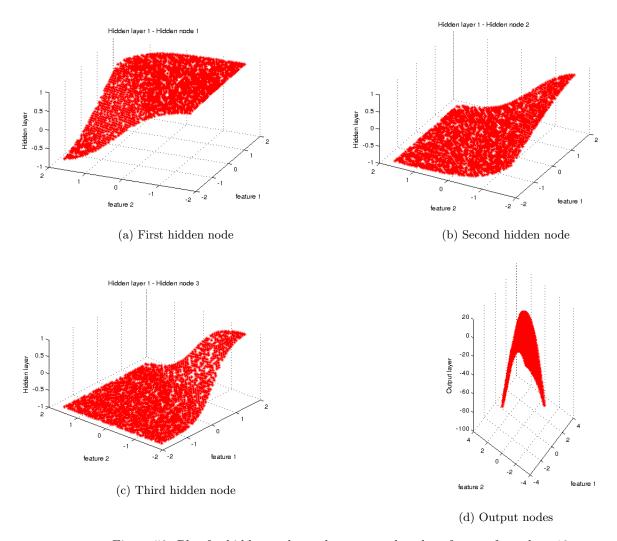


Figure 58: Plot for hidden nodes and output node values for no of epochs =10

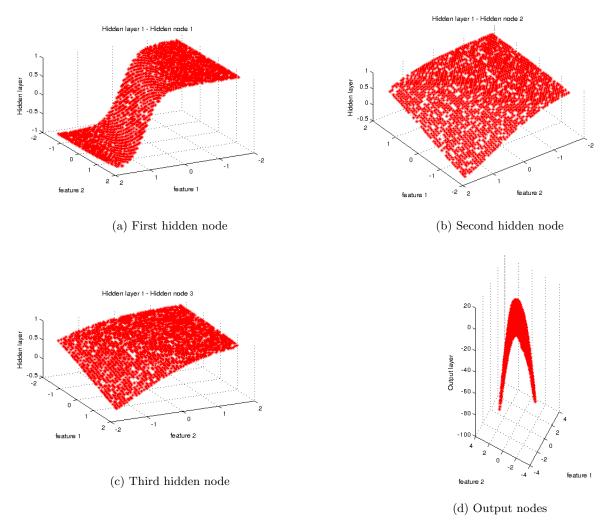


Figure 59: Plot for hidden nodes and output node values for no of epochs =50

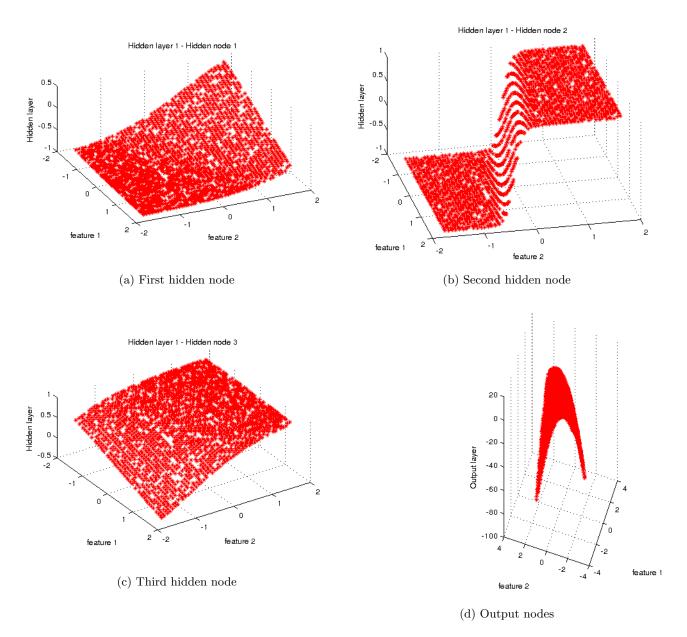
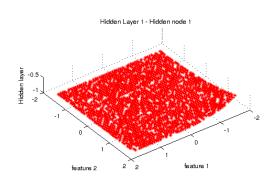
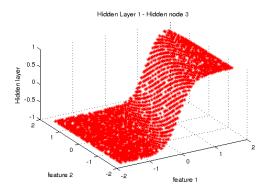


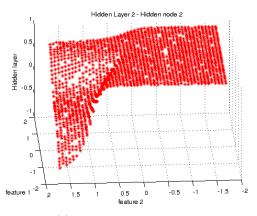
Figure 60: Plot for hidden nodes and output node values for no of epochs =100



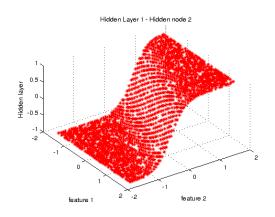
(a) Hidden layer 1 ,First hidden node



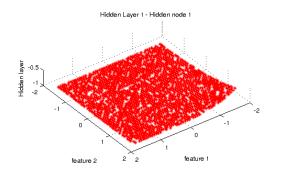
(c) Hidden layer 1, Third hidden node



(e) Hidden layer 2, Second hidden node



(b) Hidden layer 1, Second hidden node



(d) Hidden layer 2, First hidden node

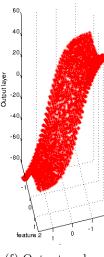


Figure 61: Plot for hidden nodes and output node values for no of epochs =1

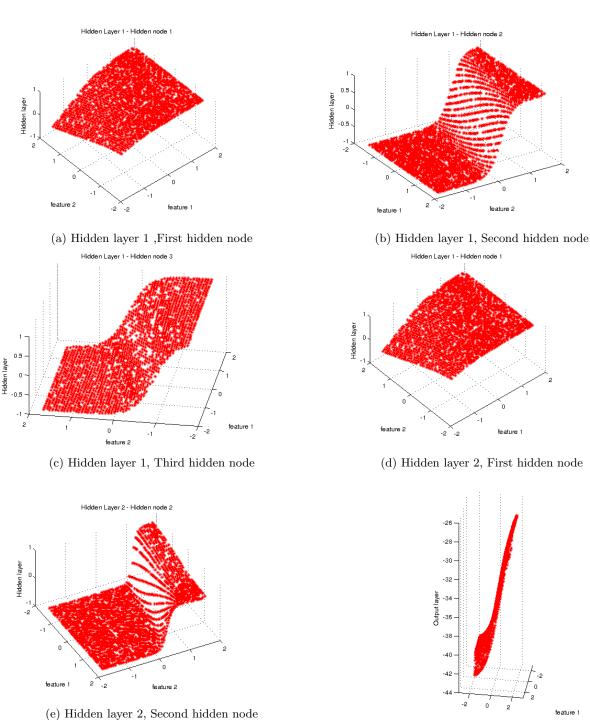
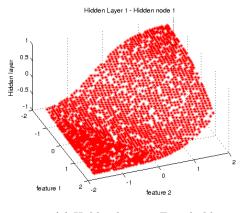
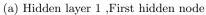
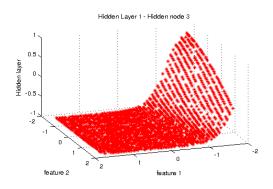


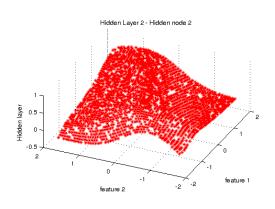
Figure 62: Plot for hidden nodes and output node values for no of epochs =2



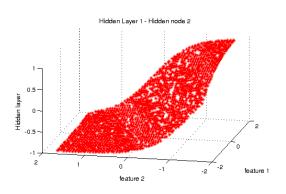




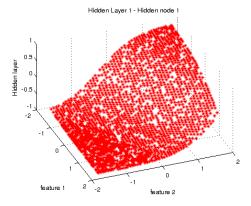
(c) Hidden layer 1, Third hidden node



(e) Hidden layer 2, Second hidden node



(b) Hidden layer 1, Second hidden node



(d) Hidden layer 2, First hidden node

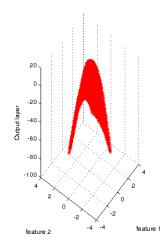


Figure 63: Plot for hidden nodes and output node values for no of epochs =10

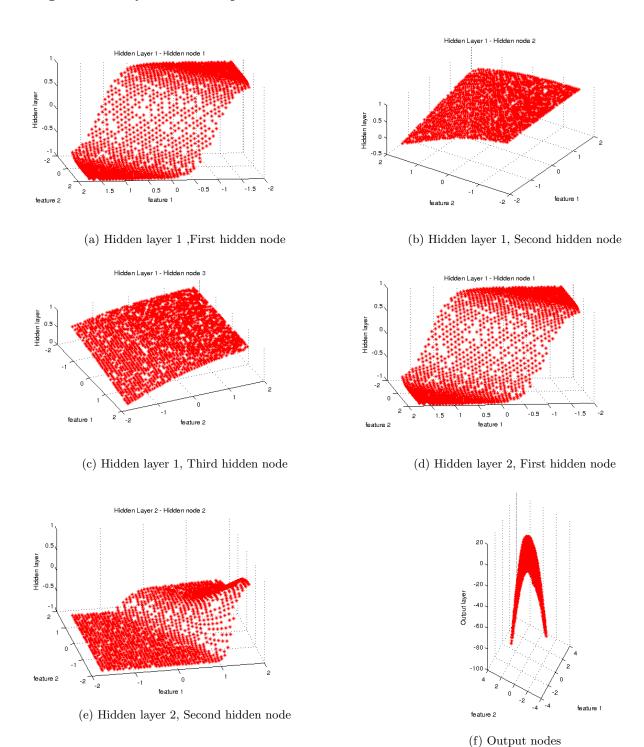
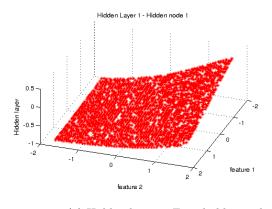
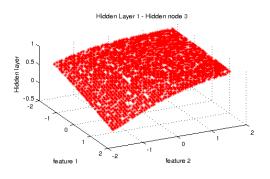


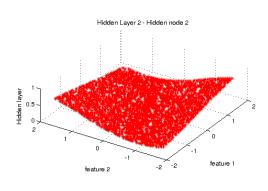
Figure 64: Plot for hidden nodes and output node values for no of epochs =50



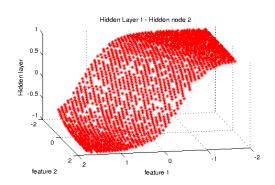
(a) Hidden layer 1 ,First hidden node



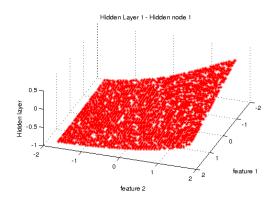
(c) Hidden layer 1, Third hidden node



(e) Hidden layer 2, Second hidden node



(b) Hidden layer 1, Second hidden node



(d) Hidden layer 2, First hidden node

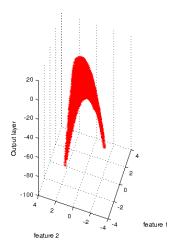


Figure 65: Plot for hidden nodes and output node values for no of epochs =100

• As the number of epochs increases, the output surface is becoming close to the underlying surface

## 3.3 Generalised RBF model

In Generalised RBF, we can use a subset of the data points itself for the RBF neurons. Selection of these RBF neurons is crucial because we can get better performance when these data points are located as diverse as possible. Also, the spread parameter should be such that interpolation can happen smoothly.

We have used K-means to find the cluster centroids for RBF neurons instead of random selection of data points. The spread is  $1/2\sqrt(M)$  times the maximum inter-centroid distance.

#### 3.3.1 Univariate Dataset

#### MSE variation with model complexities and regularisation parameter

We plotted MSE with varoius model complexities for dataset size 10 without regularization. The plot is shown below.

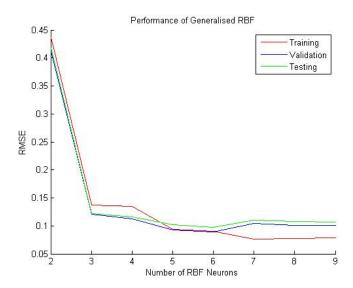


Figure 66: Performance of RBF for different numbers of RBF neurons without regularisation on N=10

#### Inferences

- Minimum error obtained for 6 neurons in the hidden layer
- We could see that overfitting has occurred at this stage and for further increase in number of RBF neurons, the training error goes down while the validation and training error are increasing

Then we plotted with regularization for  $\lambda = 0.1$ .

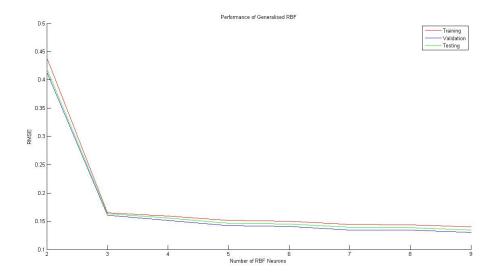


Figure 67: Performance of RBF for different numbers of RBF neurons with regularisation 0.1 on N=10

• We can see that the addition of a small regularisation term has helped to overcome the over fitting effect.

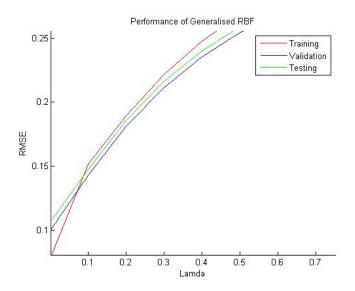


Figure 68: Performance of RBF for different values of regularisation on N=10 with 9 RBF Neurons

## Inferences

• When regularisation term is zero, the error in test and validation data are more than in training data

- When regularisation is near 0.1, the error in test data and validation data are almost same as in training data
- $\bullet$  When regularisation is crossing 0.1, the error is getting increased continuously.

# Plots of Model output Vs Target output

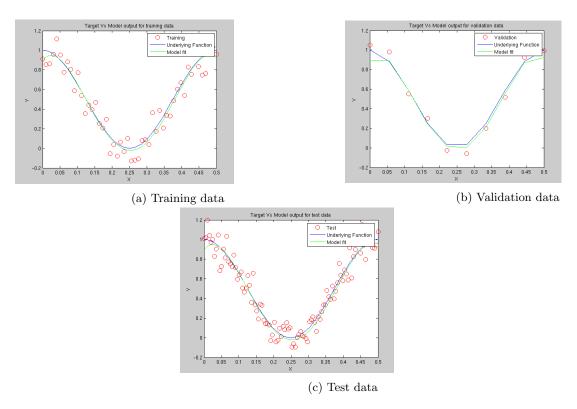


Figure 69: Model output Vs Target output for N=100 , 4 RBF Neurons

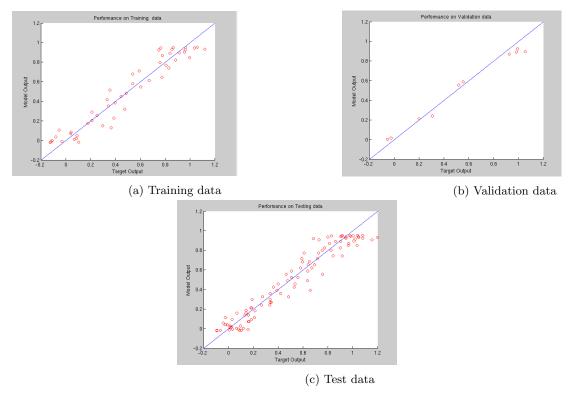


Figure 70: scatter plot with target output on x-axis and model output on y-axis for training data, validation data and test data with training data size=100

We can see that most of the scatter points are aound the x=y line showing good accuracy

#### 3.3.2 Bivariate Dataset

## MSE variation with model complexities and regularisation parameter

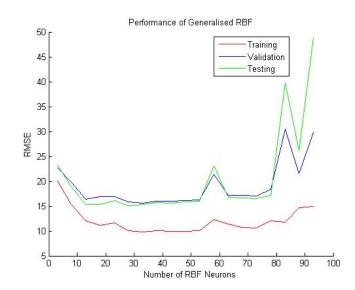


Figure 71: Performance of RBF for different numbers of RBF neurons without regularisation on N=100

#### Inferences

 $\bullet$  We can that as the model complexity, i.e. the number of RBF neurons increase in number, the RMSE on training data keeps decreasing while that on the test and validation error goes up after M=80

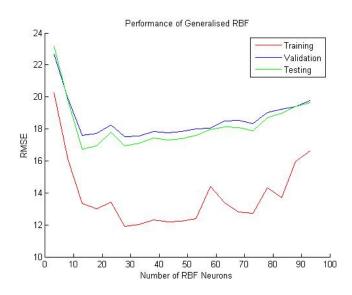


Figure 72: Performance of RBF for different numbers of RBF neurons with regularisation 0.9 on N=100

 $\bullet$  Adding a regularisation term of 0.9 on the training sample size N=100 gives better performance of the model on test data even for higher complexities

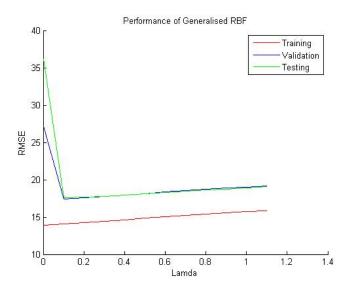


Figure 73: Performance of RBF for different values of regularisation on N=100 with 99 RBF neurons

### Inferences

- Regularisation value of 0.1 has the most effect on reducing overfitting with M=99 RBF neurons for training sample of N=100.
- With regularisation, the variance of the model has come down.

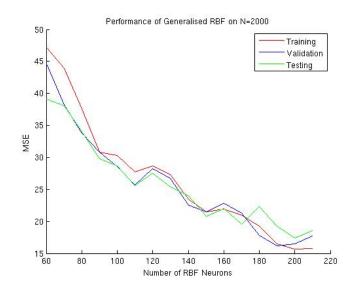


Figure 74: Performance of RBF Vs Number of RBF Neurons on N=2000

 $\bullet$  Best model performance of 16.9 MSE is achieved by choosing M=190 based on validation data performance

# Plots of Model output Vs Target output

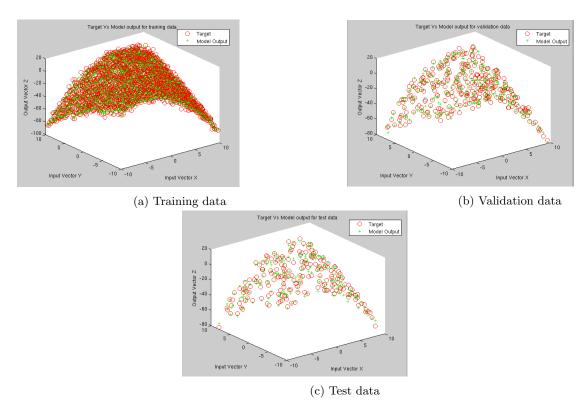


Figure 75: Model output Vs Target output for N=2000, RBF Neurons= 190

## Inferences

- Best accuracy of MSE 16.9 is obtained with 190 RBF neurons.
- This is performing better than gradient descent MLFNN on small size training sets because we are using regularisation in RBFNN.
- The newrb module in matlab could give the same accuracy with 69 neurons itself. Selection of RBF neurons in newrb module is based on picking the data point with maximum error in each iteration.

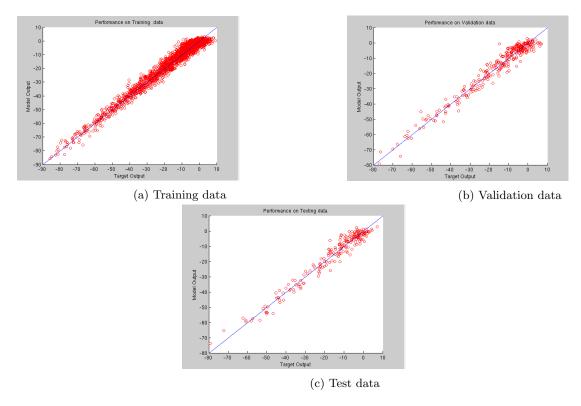


Figure 76: scatter plot with target output on x-axis and model output on y-axis for training data, validation data and test data with training data size=2000

• Most of the scatter plots are close to the x=y line.