Neural Network Machine Learning and Dimension Reduction for Data Visualization

Charles A. Liles
Langley Research Center, Hampton, Virginia
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Langley Research Center, Hampton, Virginia
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Many of the machine learning models developed for this project were created using Weka software developed by the University of Waikato. This software is available under the GNU General Public License 2.0. This project also includes software developed by the Delft University of Technology. The software developed for this project is specifically designed to support NASA’s Multi-mission Systems Analysis for Planetary Entry (M-SAPE) program.

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I. Abstract

Neural network machine learning in computer science is a continuously developing field of study. Although neural network models have been developed which can accurately predict a numeric value or nominal classification, a general purpose method for constructing neural network architecture has yet to be developed. Computer scientists are often forced to rely on a trial-and-error process of developing and improving accurate neural network models. In many cases, models are constructed from a large number of input parameters. Understanding which input parameters have the greatest impact on the prediction of the model is often difficult to surmise, especially when the number of input variables is very high. This challenge is often labeled the “curse of dimensionality” in scientific fields [1].

However, techniques exist for reducing the dimensionality of problems to just two dimensions. Once a problem’s dimensions have been mapped to two dimensions, it can be easily plotted and understood by humans [2]. The ability to visualize a multi-dimensional dataset can provide a means of identifying which input variables have the highest effect on determining a nominal or numeric output. Identifying these variables can provide a better means of training neural network models; models can be more easily and quickly trained using only input variables which appear to affect the outcome variable. The purpose of this project is to explore varying means of training neural networks and to utilize dimensional reduction for visualizing and understanding complex datasets.

Nomenclature

<table>
<thead>
<tr>
<th>BH-SNE</th>
<th>Barnes-Hut t-Distributed Stochastic Neighbor Embedding</th>
</tr>
</thead>
<tbody>
<tr>
<td>FPA</td>
<td>Flight Path Angle</td>
</tr>
<tr>
<td>LHS</td>
<td>Latin Hypercube Sampling</td>
</tr>
<tr>
<td>MLP</td>
<td>Multilayer Perceptron</td>
</tr>
<tr>
<td>MNIST</td>
<td>Mixed National Institute of Standards and Technology</td>
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<tr>
<td>M-SAPE</td>
<td>Multi-mission System Analysis for Planetary Entry</td>
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<td>MSR</td>
<td>Mars Sample Return</td>
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<td>PCA</td>
<td>Principle Component Analysis</td>
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<tr>
<td>PHP</td>
<td>PHP: Hypertext Processing</td>
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<tr>
<td>t-SNE</td>
<td>t-Distributed Stochastic Neighbor Embedding</td>
</tr>
</tbody>
</table>
II. Introduction

II.1. M-SAPE Background

This project was initiated and implemented in support of NASA’s Multi-mission System Analysis for Planetary Entry (M-SAPE). The M-SAPE project consists of a software package which can provide preliminary analysis of entry vehicle performance and viability when used to return extraterrestrial samples to the Earth’s surface. The vehicle designs on which M-SAPE performs analysis will be used for future NASA missions to safely bring scientific samples from other planetary bodies to Earth for further study. One entry vehicle for which M-SAPE is particularly designed is the Mars Sample Return (MSR) vehicle [3].

The M-SAPE software package was developed using the Python programming language, and it is intended for use in the preliminary analysis of entry vehicle performance during the early stages of design. In its current implementation, M-SAPE has 18 input variables, such as Thermal Protection System (TPS) material, vehicle radius, payload size, etc. which are specified by the user. The M-SAPE software tests the viability of user inputs and outputs over ninety nominal and numeric performance characteristics.

The M-SAPE system can be run in parallel to evaluate a queue of differing vehicle designs. However, the evaluation time of multiple models is protracted. For example, the M-SAPE software can take several days to evaluate 20,000 different entry vehicle designs even when these evaluations are conducted in parallel on a dedicated server. The need to store previously calculated M-SAPE results and to be able to interpolate values from these results was identified during the development of this project.

To address this challenge, a user-friendly M-SAPE interface for viewing previous M-SAPE results, running new M-SAPE iterations, and providing a means of interpolating stored M-SAPE information was developed. An M-SAPE interface consisting of a PHP: Hypertext Processing (PHP, a recursive acronym) web browser interface, a MySQL database back-end, and Python code for M-SAPE analysis and contour plotting was constructed. This system provided a means for users to query and view over 80,000 previously computed M-SAPE results. Users also had the ability to run their own analysis of the M-SAPE code from the PHP interface. Also, the system provided a means of contour plotting of stored M-SAPE data onto a two-dimensional plot.

The M-SAPE interface includes the ability for administrators to upload any dataset consisting of nominal and numeric data via the PHP browser interface. Datasets had to be in a pre-specified comma separated values (.CSV) file format. The PHP interface uploaded a .CSV file from the administrator’s
computer system, parsed it to determine variable names and variable types, and then built a table structure in the MySQL backend formatted to hold the uploaded data. The data in the .CSV file would then be stored into the MySQL backend. Indexes would also be automatically created on variables which had been marked as independent (or input) variables in the .CSV file. This system provided a streamlined mechanism for uploading heterogeneous datasets into a system from which data could be easily queried and plotted by a user.

II.2. Machine Learning

Machine learning tools were also integrated into the M-SAPE interface. The primary machine learning technique utilized was Multi-Layer Perceptron (MLP) neural networks developed using open-source Java-based Weka software. An MLP neural network consists of an input layer, one or more hidden layers, and an output layer [2].

![MLP Neural Network Generic Architecture](image)

**Figure 1. MLPA neural Network Generic Architecture**

The input layer is made up of individual nodes, one node for each input parameter which will be used to predict the output. The hidden layer(s) contain multilayer perceptron nodes which take input values from the previous layer, calculate an output value, and then feed the output value forward to the next layer. The output layer consists of a single node (for nominal predictions, it may contain multiple nodes, one for each potential nominal output value) which receives input values from the previous hidden layer and outputs the prediction for the model [2].

MLP neural networks must first be trained with representative data before they can be used to make prediction. Neural networks evaluate training data and adjust node weights through a means called backpropagation [2]. A Neural Network M-SAPE interface was developed which enabled administrators
to specify the number of hidden layers and nodes for a neural network design. The M-SAPE interface would then train a neural network for each numeric output variable in the dataset as well as a single nominal predictor pre-specified by the administrator. Using a two-hidden layer geometry with 25 nodes in the first layer and two nodes in the second, the M-SAPE interface took just over an hour to train 85 neural networks for the output variables belonging to the most current entry vehicle dataset uploaded into M-SAPE. The 84 numeric prediction neural networks often had a relative absolute error within 5% of the actual value during testing. The single nominal output predictor correctly predicted approximately 95% of its tested instances. Furthermore, these neural network evaluation times for prediction were much shorter than a full M-SAPE run. For example, evaluating 80,000 atmospheric entry vehicle designs took two weeks using the original M-SAPE software. However, the same evaluation would take seconds using the M-SAPE Neural Network system. These neural networks trade a relatively small amount of accuracy for a dramatic decrease in evaluation time in comparison with deterministic, M-SAPE software calculations.

The M-SAPE neural network interface enabled users to produce contour plots of user-selected variables in two dimensions. It also provided nominal shading of the contour plot to indicate design areas where entry vehicles would not converge or where TPS limits had been violated (see Figures A.1 and A.2 in the Appendix for an example of a neural network contour plot and a KD-Tree contour plot, respectively).

Although the M-SAPE interface provided a relatively efficient means of parsing, storing, querying, plotting, and predicting uploaded datasets of varying schemas, further testing was conducted to determine means of improving neural network accuracy. Testing was not just limited to the M-SAPE entry vehicle dataset. Other datasets were loaded into the M-SAPE interface for neural network training evaluation.

II.3. Dimensional Reduction

Attempts at dimensional reduction using methods such as Principle Component Analysis (PCA) and t-Distributed Stochastic Neighbor Embedding (t-SNE) were utilized in conjunction with neural network training to increase model prediction accuracy [4]. MLP neural network prediction accuracy depended greatly on neural network architecture, preprocessing of data, and the type of problem for which the network was developed.
III. Datasets

III.1. M-SAPE Entry Vehicle Dataset

As previously mentioned, M-SAPE datasets have 18 user-specified input variables. The M-SAPE software will test the viability of user inputs and output over ninety nominal and numeric performance characteristics.

The nominal outputs include vehicle convergence (whether or not the vehicle based on user input parameters will be a viable design), TPS warnings (violations of the desired TPS model), etc. The numeric outputs includes maximum total heat rate during entry, maximum entry load (measured in G-forces), total entry mass, etc. M-SAPE entry vehicle datasets are nonlinear in nature.

III.2. Vehicle Side-Impact Crashworthiness Dataset

Vehicle side impact crashworthiness datasets were also generated using a Monte Carlo analysis technique described in Ref. [5]. These crashworthiness datasets were created to analyze four-wheeled vehicles (i.e., cars), not entry vehicles. Unlike the entry vehicle data, the crashworthiness datasets were linear in nature [5].

The purpose of using the side-impact crashworthiness equations was to determine if neural network approximation could match the accuracy of the Monte Carlo prediction using the equations.

III.3. LHS Datasets

LHS is a means of producing a smaller dataset that is still representative of all input variables over either a normal or uniform distribution [6]. The goal of utilizing LHS was to minimize the number of instances required to train a neural network. If a neural network could be trained to produce accurate predictions using a smaller dataset, the time and computational requirements for training could be dramatically reduced.

Datasets for entry vehicles and vehicle side-impact crashworthiness were generated using LHS.
III.4. MNIST Hand-Drawn Digit Grayscale Dataset

The MNIST dataset is a standard machine learning dataset provided by Ref. [8]. The dataset consists of grayscale data for hand-drawn digits, and it contains 784 input variables which can be used to predict what digit was drawn. The MNIST dataset was used to evaluate neural network training on an extremely high-dimensional problem.

III.5. M-SAPE Data Upload Interface

The PHP M-SAPE interface was utilized to efficiently upload the heterogenous datasets into the M-SAPE MySQL backend. Once data had been loaded into the system, neural network models could be expediently trained and evaluated for prediction accuracy.

IV. Methods

IV.1. Neural Networks

Weka software was used for the training and evaluation of MLP neural networks. Training of these networks was directly integrated into the M-SAPE PHP interface. Trained models could then be stored on the M-SAPE server and utilized for predictions and contour plotting. The Weka software utilized was capable of producing both nominal and numeric value predictions. Training with standard entry vehicle datasets and with LHS-generated datasets was conducted.

IV.2. Data Pre-Processing and Dimensional Reduction

Techniques were also incorporated for pre-processing datasets prior to neural network training. PCA and removal of non-essential variables were just two means of refining training datasets prior to MLP network training with the goal of improving neural network accuracy.

Also, t-SNE was utilized for reducing multi-dimensional problems to just two dimensions for plotting and data visualization [4]. Although an effective means for dimension reduction, t-SNE was computationally cumbersome (running in O(n^2) time). A more efficient version of t-SNE known as Barnes-Hut t-Distributed Stochastic Neighbor Embedding (BH-SNE) was utilized in lieu of standard t-
SNE. This algorithm runs in O(n log n) time and is an effective means of quick dimension reduction with little loss in fidelity compared with standard t-SNE [7].

Both t-SNE and BH-SNE are unsupervised methods of dimension reduction. Unlike a neural network which compares a prediction with the actual value and then adjusts each node using backpropagation, the SNE algorithms simply group input instances together in two dimensional space based on the values of their input variables. The actual output value is never referenced during this process of dimension reduction. However, the results can then be plotted in a two dimensional scatter plot with a nominal color coding to show the actual output. If similar colors are clustered together, the algorithm has effectively mapped a multi-dimensional problem into just two dimensions.

SNE techniques have often been used for unsupervised learning of gray scale images (such as the MNIST dataset). For example, previous studies have shown t-SNE’s effectiveness at reducing over a hundred dimensions representing grayscale, hand-written digits to just two dimensions. The results revealed ten clusters representing each digit where like digits are clustered together [4].

During this set of tests, BH-SNE was used to not only map a nominal output for entry vehicle performance, but to also map input variable values to separate scatter plots with the same plot pattern as the output scatter plot. By comparing the output scatter plot with each input scatter plot, a user could easily discover which input variable appears to have the highest effect on the output variable. A similar colored scatter pattern between an output variable and a particular input variable would indicate a strong relationship between the output and the input variable. This technique could be particularly useful in multi-dimensional nonlinear problems where identifying which input variable has the highest effect on the output is especially difficult.

V. Results

The ability of a neural network to provide accurate predictions depended greatly on the type of problem being predicted, pre-processing techniques used prior to training, utilization of representative sampling for network training, and the network architecture. Although PCA proved to be an effective pre-processing technique for linear problems like vehicle side-impact crashworthiness, it actually decreased prediction accuracy for nonlinear problems such as the M-SAPE entry vehicle dataset. Improvements in nonlinear problem models could only be achieved by varying the architecture of the neural network. While PCA was invaluable for identifying which input variables had the greatest impact on the output for a linear problem, BH-SNE proved to be more effective for identifying relationships between input and output in nonlinear problems. Using LHS proved to be an effective means of reducing the required number of instances needed for MLP network training.
V.1. Neural Network Testing Results for Linear Problems

As mentioned above, multiple vehicle side impact crashworthiness datasets were used to build neural networks. These datasets contained nine differing numeric input variables that determine 11 different numeric outputs. However, for each output variable, only a subset of the input variables were actually used to calculate the output (using the Monte Carlo analysis presented in Ref. [5]). To determine the accuracy of neural network prediction, all input variables were initially used to produce models for each output variable. If the neural network backpropagation method was effective, then input variables should be appropriately weighted according to their relationship to the output variable.

To test neural network prediction with datasets of differing sizes, three separate neural networks were trained: one with only 10 data points, one with 100, and another with 1,000 data points. These datasets were created using LHS over a uniform distribution of the input variables. A dataset of 10,000 instances was then used to test each model. In theory, if more data points are used to train a neural network model, the model should be more accurate. Figure 2 is a plot for each models’ resulting mean and standard deviation for Load (Abdomen) prediction (in red) compared with the Monte Carlo analysis actual expected value (in blue).

![Figure 2. Comparison of Mean and Standard Deviation for Neural Networks Trained with Varying Numbers of Instances (X-Axes are on a logarithmic scale)](image-url)
As the number of data points used to train each neural network increases, the mean and standard deviation for its predictions more closely match those values produced by the Monte Carlo formulas used to compute the actual values (shown in blue). However, it was possible to use LHS to reduce the number of instances required for training (with a trade off in prediction accuracy).

As mentioned above, not all input variables were used to calculate certain outputs. For example, the Load (Abdomen) function was as follows:

\[
\text{Load (Abdomen)} = 1.16 - 0.3717 \ d_2 \ d_4 - 0.00931 \ d_2 \ d_{10} - 0.484 \ d_3 \ d_9 + 0.01343 \ d_6 \ d_{10}
\]

For this calculation, only the \(d_2, d_4, d_{10}, d_3, d_9, \) and \(d_6\) variables were used to produce the actual result using Monte Carlo Analysis. However, the neural networks trained for this example utilized every variable for training of the models. If the backpropagation weighting method worked correctly during training, then the neural network should have appropriately compensated for input variables which have no impact on the actual output.

Figure 3 includes contour plots produced by the 10-instance training set model, the 100-instance training set model, and the 1,000-instance training set model created in M-SAPE for Load (Abdomen) predictions:

![Figure 3. Load (Abdomen) Crashworthiness Variable Contour Plots for Neural Networks Trained with Varying Numbers of Instances](image)

For the contour plots in Figure 3, \(d_2\) is the Y-Axis and \(d_1\) is the X-Axis. All other input variables are set at their mean values according to Ref. [5]. We know that \(d_2\) is used to calculate Load (Abdomen), but \(d_1\) is not. So, if the neural network backpropagation training method is effective, the slope of the contour lines should be completely flat (because \(d_1\) should have no impact on the prediction). As the number of instances used to train the neural network increases, the accuracy of the contour plots also increases. In the 10-instance trained model, the contour lines have a definite downward slope. For the
100-instance model, the lines have a very slight upward slope. The 1,000-instance trained model has an even smaller upward slope than the 100-instance neural network. As expected, the neural network backpropagation training technique is more effective when a larger dataset is used for training.

As mentioned above, PCA testing was also used in conjunction with neural net training. The PCA technique pre-weights input variables based on their relationship to the output. Although PCA may not completely eliminate an unnecessary input variable, it will give that variable an appropriately small weight prior to feeding it to a neural network for training [2].

To test the effectiveness of PCA in conjunction with neural network training, four separate tests were conducted to predict Cost(Weight) (as defined in Ref. [5]) using the following criteria:

1. 1,000 data points, no PCA, all input variables
2. 1,000 data points, PCA, all input variables
3. 1,000 data points, no PCA, only significant variables (only those variables which are actually used in the Monte Carlo function to produce the output)
4. 1,000 data points, PCA, only significant variables

A plot of the mean and standard deviation of the results compared with the original Monte Carlo analysis is provided in Figure 4 below.

![Figure 4. Cost(Weight) Mean and Standard Deviation for Neural Networks with and without PCA and with and without Significant Variables versus Monte Carlo Prediction](image-url)
The horizontally adjacent plots in Figure 4 show the same data. The left graph uses the same logarithmic X-axis as used in Figure 2 above, and the right graph shows an enlarged version of the mean and standard deviation for the 9,995th testing data point upwards. On the left graph, the neural networks trained using PCA with all variables, no PCA with only significant variables, and PCA with only significant variables are plotted directly on top the Monte Carlo Analysis information (the enlarged graph on the right is displayed to show the distinction between these four plots). The neural network with PCA using all input variables most closely approximates the expected Monte Carlo actual value for mean values. This value differs by less than 0.02% of the Monte Carlo analysis mean for the 9,999th instance (reference the mean graph on the right above). The neural network without PCA trained only with significant variables most closely approximates the standard deviation of the Monte Carlo analysis. This neural network’s standard deviation prediction had only a difference of ~0.5% from the Monte Carlo analysis standard deviation for the 9,999th instance. Figure 5 below illustrates a plot of the mean absolute error for all models [2].

![Mean Absolute Error](image)

**Figure 5. Cost(Weight) Mean Absolute Error for Neural Networks with and without PCA and with and without Significant Variables**

The two plots represent the same data; the Y-axis in the graph on the right has been compressed. The X-axis is logarithmic for both graphs. Mean absolute error is calculated as follows:
\[
\sum_{i=1}^{n} \left| p_i - a_i \right| \over n
\]

Where \( p \) represents the predicted value, \( a \) represents the actual value, and \( n \) is the total number of evaluated instances. This form of error evaluation diminishes the effect of outlier data on the overall evaluation of models [2]. In the plots above, the model trained using all input variables with PCA applied first outperforms the other models. This mimics the results of the plots of mean in Figure 4.

It should be noted that all of the neural networks trained for this test performed relatively well. The neural network not using PCA with all input variables consistently provided the least accurate results. However, this level of inaccuracy had a difference in mean values from the Monte Carlo results of less than 1% and a difference in standard deviation of less than 10%.

In the results illustrated in Figure 4, it was known before training which variables were significant and which were not. Therefore, testing could be conducted with and without insignificant variables to determine the MLP neural network’s ability to identify and weight extraneous inputs appropriately. Surprisingly, the neural network trained with all input variables and with PCA pre-processing provided mean values closest to the expected Monte Carlo prediction. As for the standard deviation prediction, it was only slightly outperformed by the model built using only significant variables without PCA pre-processing. PCA offered an excellent means of pre-weighting input variables for better neural network training for linear problems.

V.2. Neural Network Testing Results for Nonlinear Problems

Entry vehicle datasets produced by M-SAPE were utilized as a nonlinear problem dataset for neural network training and testing. This dataset is fundamentally different than the vehicle side impact Monte Carlo analysis model used for linear problem testing. The entry vehicle dataset included 18 different input variables of both numeric and nominal data types. The numeric input variables included FPA, entry velocity, input vehicle diameter, payload mass, etc. The nominal input variables consisted of the TPS material used for both the fore and aft body of the entry vehicle and whether the craft was designed to meet MSR requirements.

As mentioned above, the M-SAPE software designed to predict entry vehicle performance will output approximately a hundred numeric and nominal variables describing the craft’s characteristics. Examples of numeric output variables are forebody maximum total heat rate, maximum entry load, and total entry mass. The nominal variables include vehicle convergence, which states if the vehicle defined
per the user’s parameters will actually be a valid model. Other nominal output variables include forebody and aftbody TPS warnings, which are textual descriptions of any potential TPS warnings for a vehicle on entering the Earth’s atmosphere.

The deterministic M-SAPE vehicle prediction software iteratively adjusted a vehicle’s input parameters if the initial specification did not produce a converged model. The process of evaluating a single model took several minutes to complete. Due to the complexity of producing the output of a vehicle’s performance, the M-SAPE entry vehicle dataset was excellent for the training and testing of neural networks for nonlinear problems.

As noted above, PCA is a useful technique for pre-processing datasets prior to neural network training for linear problems. To evaluate its effectiveness for a nonlinear problem, two separate MLP networks were built to predict forebody maximum total heat rate. One model had PCA pre-processing and the other did not. Both neural networks had a two-hidden layer architecture with 25 nodes in the first layer and two in the second. Table 1 below illustrates the difference in error between the models during testing post-training (a 67% / 33% training / testing split was used for all neural network models designed for numeric prediction) [2].

<table>
<thead>
<tr>
<th>Neural Network with PCA</th>
<th>Neural Network without PCA</th>
</tr>
</thead>
<tbody>
<tr>
<td>Correlation coefficient</td>
<td>0.9972</td>
</tr>
<tr>
<td>Mean absolute error</td>
<td>21.2896</td>
</tr>
<tr>
<td>Root mean squared error</td>
<td>36.1551</td>
</tr>
<tr>
<td>Relative absolute error</td>
<td>6.7241 %</td>
</tr>
<tr>
<td>Root relative squared error</td>
<td>8.6551 %</td>
</tr>
<tr>
<td>Total Number of Instances</td>
<td>7258</td>
</tr>
</tbody>
</table>

**Table 1. Error during Training for Neural Networks with and without PCA Pre-Processing**

Unlike the linear problem described in subsection V.1, the neural network trained without using PCA performed much better than its PCA counterpart. In another test, both the PCA and non-PCA models were evaluated with 21,348 test instances. A comparison of the mean absolute error is illustrated in Figure 6 below [2].
In Figure 6, the Y-axis represents mean absolute error while the X-axis is the number of instances evaluated on a logarithmic scale. For this example, the neural network trained \textit{without} using PCA preprocessing consistently has a lower mean absolute error than the PCA model.

LHS was also used to reduce the number of instances needed to train a neural network for a nonlinear problem. As of February 2014, the M-SAPE prediction software took a uniform distribution of all numeric inputs. New prediction datasets were produced using the Cartesian product of all perturbations in input variables. This process created a dataset of over 80,000 individual vehicle entry predictions. However, in the future, more adjustable input variables will be added to the M-SAPE software. Using a Cartesian product of perturbations in an increasing number of input variables will create massive datasets. These massive datasets will require more space for storage and will require much more computational power initially to produce and then to use for the training of neural networks. On the other hand, if LHS could be used to maintain a relatively low number of entry vehicle instances while adding new input variables, the storage and computational requirements for building neural networks could be greatly reduced.

Two datasets were created utilizing LHS with 100 and 1,000 entry vehicle designs. Neural networks were then trained from both LHS datasets as well as the previous Cartesian product dataset originally implemented with M-SAPE to predict max entry load (measured in G-forces). A separate testing dataset consisting of 5,040 entry vehicle instances was calculated using the M-SAPE
deterministic software. All neural network models were tested against the 5,040 test dataset. The mean absolute error calculations from these tests are presented in Table 2 below.

<table>
<thead>
<tr>
<th></th>
<th>100 Instance LHS Dataset Neural Network</th>
<th>1000 Instance LHS Dataset Neural Network</th>
<th>Cartesian Product Dataset Neural Network</th>
</tr>
</thead>
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<tr>
<td>Correlation coefficient</td>
<td>0.954</td>
<td>0.9978</td>
<td>0.9998</td>
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<tr>
<td>Mean absolute error</td>
<td>7.4034</td>
<td>1.1751</td>
<td>0.4622</td>
</tr>
<tr>
<td>Root mean squared error</td>
<td>10.8238</td>
<td>2.3786</td>
<td>0.7135</td>
</tr>
<tr>
<td>Total Number of Instances</td>
<td>5040</td>
<td>5040</td>
<td>5040</td>
</tr>
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</table>

Table 2. Comparison of Neural Networks Trained with LHS Datasets and Cartesian Product Datasets (Neural Networks Trained to Predict Max Entry Load (G’s))

For the 5,040 instance dataset used in the above predictions, the average actual value for all instances was 71.039 G’s. For the example above, the mean absolute error and root mean squared error (another means of measuring prediction accuracy described in Ref. [2]) were smallest when using the neural network trained from the large, Cartesian product dataset of 5,040 instances. However, the MLP network trained from the 1,000 instance LHS dataset provided fairly accurate predictions compared to its Cartesian product counterpart. The 1,000 instance neural network used a LHS generated dataset 1/5th the size of the Cartesian product dataset, but it was fairly accurate. On the other hand, using the neural network trained from the 100 instance LHS dataset was much less accurate. For the 100 instance dataset, the mean absolute error was just over 10 % of the average value in the test dataset.

Figure 7 below illustrates a plot of the mean absolute error values from the tests described above.

Figure 7. Comparison of Mean Absolute Error for Neural Networks Trained with LHS Datasets and Cartesian Product Datasets (Neural Networks Trained to Predict Max Entry Load (G’s))
Although the neural network trained with 5,040 instances provided the most accurate results, the 1,000 instance LHS dataset provided a fairly small loss of accuracy while using much less data to train the MLP network model. The network trained with the 100 instance LHS dataset, on the other hand, produced a much higher deviation in mean absolute error than the other two networks.

V.3. Neural Network Testing Results for Nonlinear, High Dimensional Problems

Testing on nonlinear datasets of high dimensionality were also conducted using the MNIST dataset. This dataset had 784 input variables used to predict a hand-written digit [8]. Using a neural network consisting of two layers with 25 nodes in the first layer and two nodes in the second, a neural network model with an accuracy of 76.4706% was created. Another two layer model with 50 nodes in the first layer and four nodes in the second layer accurately predicted the drawn digit for 79.5294% of the test instances. Yet another two layer model with 100 nodes in the first layer and eight nodes in the second layer provided a prediction accuracy of 87.1765%. Varying the network geometry by increasing the number of nodes in each layer increased the accuracy of models built for extremely high dimensional problems.

V.4. Nonlinear Dimension Reduction and Visualization

Although PCA failed to produce more accurate prediction models for nonlinear problems, BH-SNE was also utilized to reduce multidimensional problems to just two dimensions for neural network training as well as for plotting. As mentioned above, BH-SNE is an unsupervised method of reducing dimensionality. BH-SNE is also specifically implemented for nominal value predictions. [7]

To evaluate the effectiveness of BH-SNE dimension reductions, a dataset for predicting entry vehicle convergence and TPS warning messages was produced using BH-SNE dimensional reduction to create just two input variables. Two nominal neural network prediction models were then created, one using the BH-SNE dataset and the other using all original input variables. A two--layer architectures of 25 nodes in the first layer and two nodes in the second were implemented for both models. Table 3 below shows a comparison of the results for training both models.
Clearly, the neural network trained without any BH-SNE preprocessing performed much better than its counterpart with BH-SNE dimensional reduction. The network trained without BH-SNE predicted 95.1665 % of the test instances correctly, while the network trained with BH-SNE preprocessing had a 58.5377 % prediction accuracy.

Varying network architecture did not significantly improve BH-SNE preprocessing neural network prediction. A two-layer architecture of 50 nodes in the first layer and four nodes in the second layer produced a prediction accuracy of 59.9806 %; only a small increase in accuracy from the model shown in Table 3 above. However, another two-layer architecture of 100 nodes in the first layer and four nodes in the second layer yielded a model with just a 58.6897 % prediction accuracy.

Although BH-SNE preprocessing did not have a positive effect on neural network predictions, it proved to be a valuable technique for data visualization. As mentioned above, BH-SNE reduces a multidimensional dataset to just 2 dimensions. The data can then be plotted easily on a scatter plot. BH-SNE scatter plots have an arbitrary scale for both the X and Y axes. It does not necessarily matter exactly where on the X and Y axes an instance is plotted. However, if BH-SNE scatter plotting is performed on a dataset where input values are causal to the value of the output, instances with similar input values will be clustered together on the plot [7].

Previous implementations of BH-SNE have simply plotted the output variable with specific color codes on a scatter plot. When like colors are grouped together, the BH-SNE dimension reduction process has worked successfully. For the M-SAPE BH-SNE implementation, the output variable was mapped to a scatter plot as in [7]. However, input variables were then also mapped to the same plot cluster pattern. Figure A.3 in the Appendix illustrates a BH-SNE plot of an output variable for the convergence and any TPS Warnings for entry vehicles. Figure A.4 through A.11 in the Appendix are BH-SNE scatter plots where the original input variables have also been mapped to the same plot as the output variable in Figure A.3. Reducing dimensions and then plotting the inputs and output to the same scatter plot can be used to deduce which variables have the greatest effect on the output variable.
An example of the benefits of this comparison process is provided below. This comparison focuses on the same, isolated cluster pattern across both input and output scatter plots. By comparing similarities in color patterns between the input and output plots, a user can more easily determine what variables have the greatest influence on the output variable: convergence and TPS warnings. Figure 8 below shows an initial comparison in subsections of the plots provided in Figures A.3 through A.11 in the Appendix.

![Comparison of variables](image)

**Figure 8. Comparison of Convergence and TPS Warnings, Nose Radius / Input Vehicle Radius, Aft and Forebody TPS Materials, and MSR Mode BH-SNE Scatter Plots**

Figure 8 illustrates the output variable in the top left with all other plots being input variables. By viewing the cluster patterns side by side, a user can easily determine that this cluster is for a vehicle with a .75 Nose Radius / Input Vehicle Radius ratio and an aftbody material of SLA-561V with a PICA-AL-5056 forebody. The vehicle is also designed for MSR mode.

Although the appraisal of the plots in Figure 8 elucidates the design parameters for the set of vehicles plotted for this cluster, it does not reveal what variable is specifically causing the vehicle designs to converge or not converge with or without TPS warnings. Another comparison is provided in Figure 9 to help determine what input variable is causal to the convergence of this particular type of vehicle.
By carefully comparing the two images in Figure 9 side by side, it appears that as the entry velocity of the vehicle is increased past a value of 14,000 m/s, this type of entry vehicle will still converge, but it will have a TPS warning for the forebody. The convergence with forebody TPS warnings is plotted in gold on the left. The user seeks to design a vehicle that will converge with no warnings (areas plotted in red on the left). So, for this particular entry vehicle design, the entry velocity must be 14,000 m/s or slower to have any chance of converging with no TPS warnings.

Although velocity appears to definitely play a role in forebody TPS performance in Figure 9, it still does not account for the areas of non-convergence with aft and forebody TPS warnings plotted in green on the left image. The images in Figure 10 below illustrate yet another variable that appears to directly impact this non-convergence.
In Figure 10, it appears that whenever the entry vehicle’s diameter is 2 m (plotted in green on the right), the vehicle will not converge, and it will have both fore and aft TPS warnings (also plotted in green on the left). Figure 11 below shows scatter plots of convergence and TPS warnings, vehicle entry velocity, and vehicle diameter.

**Figure 11. Comparison of Convergence and TPS Warnings, Vehicle Entry Velocity, and Input Vehicle Diameter BH-SNE Scatter Plots**

From viewing the three images side by side in Figure 11, a user can reasonably hypothesize that for this entry vehicle design, the entry velocity must be 14,000 m/s or slower and the vehicle diameter must be narrower than 2 m for the vehicle to converge with no warnings. Using the M-SAPE neural network contour plotting interface, a scatter plot was created to test the efficacy of this hypothesis (Figure 12 below).

**Figure 12. M-SAPE Neural Network Contour Plot for an MSR Mode Entry Vehicle with an SLA-561V Aftbody and PICA-AL-5056 Forebody**
In Figure 12, entry vehicle velocity is the input variable selected for the X-Axis while the Y-Axis represents input vehicle diameter. Areas shaded in gold represent vehicles that converge with a forebody TPS warning. Areas in red represent designs that will not converge; unshaded areas are vehicles that converge with no TPS warnings. A user can clearly see that as entry vehicle velocity reaches 15,000 m/s or higher, the vehicle will begin to have forebody TPS warnings (although it will still converge). As the vehicle’s diameter increases above 1.8 m, the vehicle will either have forebody TPS warnings, or it will not converge at all. The contour plot in Figure 12 appears to match the hypothesis surmised from comparing BH-SNE plots in Figure 11.

VI. Summary

The predictive success of any machine learning model will depend greatly on the type of problem for which it has been designed, data preprocessing techniques, representative sampling for the training data, and the architecture of the network. Whereas a preprocessing technique such as PCA can prove useful for a neural network trained to predict a linear problem, it often negatively impacts results for a nonlinear predictive model. LHS is an excellent means of developing a relatively smaller dataset for training that is still representative of the problem, whether it is linear or nonlinear. While BH-SNE should not be used for dimension reduction prior to neural network training, it is an excellent technique for nonlinear data visualization and data discovery.

If the problem for which a neural network is being designed is known to be linear in nature, PCA is a valid preprocessing technique which can improve predictive results. The experiments described in the first part of this technical memorandum utilized deterministic, linear equations from Ref. [5]. As illustrated in Figure 4, PCA preprocessed data often provided more accurate predictive results than datasets which had not been processed with PCA prior to training and testing.

Unfortunately, it is not always known beforehand whether the problem for which a neural network is being designed is linear or nonlinear in nature. Often, MLP networks are trained from a dataset which was produced by recorded observations, not one derived from deterministic equations such as the ones used in this memorandum. If the linearity of the problem for which the model being designed is unknown, neural network training can be conducted with and without PCA preprocessing. If a MLP network trained from PCA-processed data appears to be more accurate than its non-PCA counterpart, the problem may be linear in nature in which case PCA may be helpful for prediction accuracy. However, if PCA pre-processing does not appear to improve the prediction accuracy of the network model, the
problem may be nonlinear. For nonlinear problems, no preprocessing should be implemented for the input dataset.

For reducing the amount of data required to train a neural network, LHS is an excellent technique for either linear or nonlinear problems. A LHS dataset may be used to more easily train a neural network with only a small loss of predictive accuracy. Although the LHS sampling techniques described in this memorandum were created from a uniform distribution, a Gaussian distribution may also be utilized. While LHS was tested in writing this memorandum with the desire to minimize the computation of complex entry vehicle datasets, it could also be used for efficient design of experiments. If the number of experiments required for creating a representative dataset is minimized with LHS, the time and cost of conducting a number of experiments can be dramatically reduced [6].

While BH-SNE did not prove to be a beneficial alternative to PCA for nonlinear data preprocessing, it is an effective technique for data visualization. By mapping both input variables and output variables to the same scatter plot pattern using BH-SNE, correlation or causation between inputs and outputs can be easily determined by comparing plots (a JavaScript implementation was also created for this comparison process and can be viewed in Figure A.11 in the Appendix). Whereas the BH-SNE experiments conducted for this memorandum could be utilized to find causal relationships between input variables and the output, this may not always be the case for other problems. For example, if a dataset is built from simple observation between output and presumed inputs, similarities in BH-SNE input and output scatter plot patterns may only illustrate correlation between inputs and output. In conducting experiments for this technical memorandum, the causal nature between inputs and outputs was already known because the datasets were generated deterministically. This may not always be the case for other datasets.

Although BH-SNE did not provide a beneficial means of dimension reduction prior to neural network training, high dimensionality is not necessarily preclusive to training an effective MLP network model. For example, the MNIST dataset for hand-drawn digits from Ref. [8] contains 784 inputs. A larger degree of dimensionality was accommodated for by varying the architecture of the designed neural network. Although the time required to train a network for a high dimensional problem can be protracted, the time required to use a trained network for predicting multiple instances is negligible.

VII. Conclusions

The field of neural networks is still relatively new; there is no definitively correct means of building a neural network. While LHS and occasionally PCA may be used to decrease training time and increase accuracy, experimentation with network architecture is often the best means of improving
predictive results. The BH-SNE data visualization technique, used in conjunction with neural network testing, is also an excellent means of machine learning data discovery. Experimentation with the different techniques described in this memorandum can help improve neural network training and can illuminate causal relationships between neural network input and output variables.
Appendix

Figure A.1. M-SAPE Neural Network Contour Plot

Figure A.2. M-SAPE KD-Tree Contour Plot
Figure A.3. BH-SNE Plot for Convergence and TPS Warnings (Nominal Output Variable)

Figure A.4. BH-SNE Plot for Aft and Forebody TPS Materials (Nominal Input Variables)
Figure A.5. BH-SNE Plot for MSR and Non-MSR Modes (Nominal Input Variables)

Figure A.6. BH-SNE Plot for Flight Path Angle (Numeric Input Variable Measured in Degrees)
Figure A.7. BH-SNE Plot for Entry Velocity (Numeric Input Variable in m/s)

Figure A.8. BH-SNE Plot for Input Vehicle Diameter (Numeric Input Variable in meters)
Figure A.9. BH-SNE Plot for Nose Radius to Input Vehicle Radius Ratio (Numeric Input Variable)

Figure A.10. BH-SNE Plot Payload Diameter to Vehicle Diameter Ratio (Numeric Input Variable)
Figure A.11. BH-SNE Plot Payload Mass (Numeric Input Variable Measured in kg)

Figure A.12. JavaScript Browser Interface for BH-SNE Comparison
References


Neural network machine learning in computer science is a continuously developing field of study. Although neural network models have been developed which can accurately predict a numeric value or nominal classification, a general purpose method for constructing neural network architecture has yet to be developed. Computer scientists are often forced to rely on a trial-and-error process of developing and improving accurate neural network models. In many cases, models are constructed from a large number of input parameters. Understanding which input parameters have the greatest impact on the prediction of the model is often difficult to surmise, especially when the number of input variables is very high. This challenge is often labeled the “curse of dimensionality” in scientific fields. However, techniques exist for reducing the dimensionality of problems to just two dimensions. Once a problem’s dimensions have been mapped to two dimensions, it can be easily plotted and understood by humans. The ability to visualize a multi-dimensional dataset can provide a means of identifying which input variables have the highest effect on determining a nominal or numeric output. Identifying these variables can provide a better means of training neural network models; models can be more easily and quickly trained using only input variables which appear to affect the outcome variable. The purpose of this project is to explore varying means of training neural networks and to utilize dimensional reduction for visualizing and understanding complex datasets.

Dimension reduction; Machine learning; Neural network; Systems analysis