Machine Learning Lecture Notes

Lecture Notes by Şükrü Ozan, Ph.D.

LECTURE 08

_

Motivations Non-linear Hypotheses

Non-linear Classification $x_1 = \text{size}$ $x_2 = \#$ bedrooms

 $x_3 = \#$ floors

 $x_4 = age$

 x_{100}

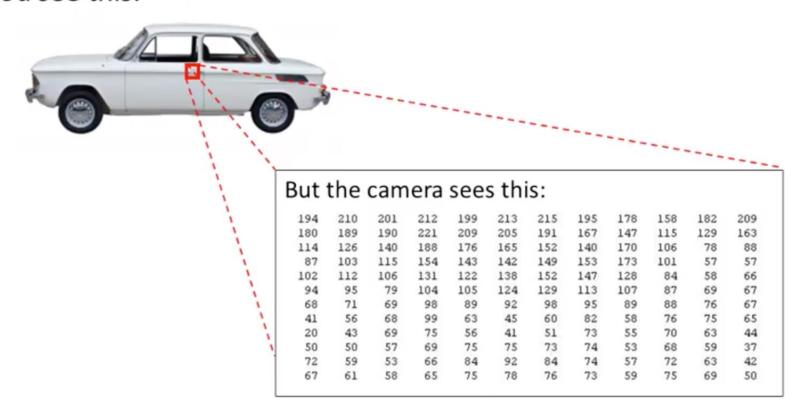
$$+\theta_{3}x_{1}x_{2}+\theta_{4}x_{1}^{2}x_{2} \\ +\theta_{5}x_{1}^{3}x_{2}+\theta_{6}x_{1}x_{2}^{2}+\dots)$$

$$\times_{1},\dots,\times_{100},\times_{1}^{2},\dots,\times_{100},\times_{1}^{2},\dots,\times_{100},\times_{1}^{2},\dots,\times_{100},\times_{$$

 $g(\theta_0 + \theta_1 x_1 + \theta_2 x_2)$

What is this?

You see this:



Computer Vision: Car detection

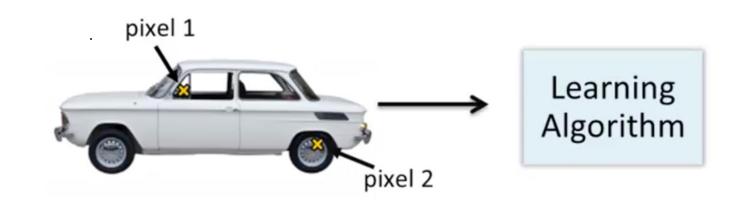




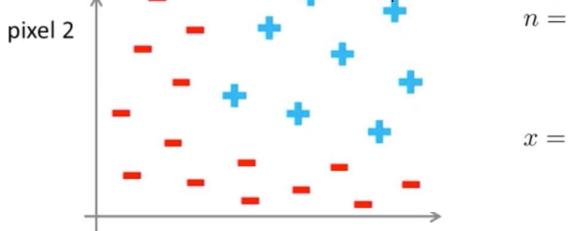
Testing:



What is this?



pixel 1



Cars

"Non" Care

n=2500 (7500 if RGB)

50 x 50 pixel images \rightarrow 2500 pixels

Quadratic features ($x_i \times x_j$): \approx 3 million features

Suppose you are learning to recognize cars from 100×100 pixel images (grayscale, not RGB). Let the features be pixel intensity values. If you train logistic regression including all the quadratic terms ($x_i x_j$) as features, about how many features will you have?

- O 5,000
- 0 100,000
- \bullet 50 million (5 imes 10^7)
- \bigcirc 5 billion (5 imes 10^9)

_

Motivations Neurons and Brain

Neural Networks

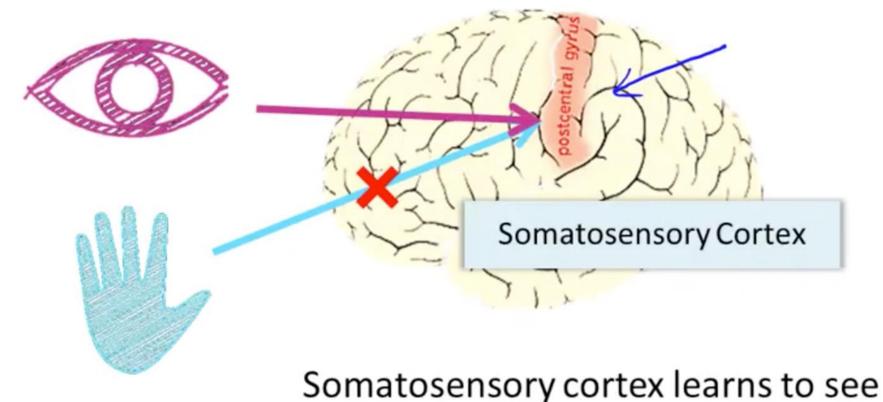
Origins: Algorithms that try to mimic the brain.

Was very widely used in 80s and early 90s; popularity

diminished in late 90s.

Recent resurgence: State-of-the-art technique for many applications

The "one learning algorithm" hypothesis



[Rolectine 8a F. r 69921 989]

Sensor representations in the brain





Seeing with your tongue





Haptic belt: Direction sense



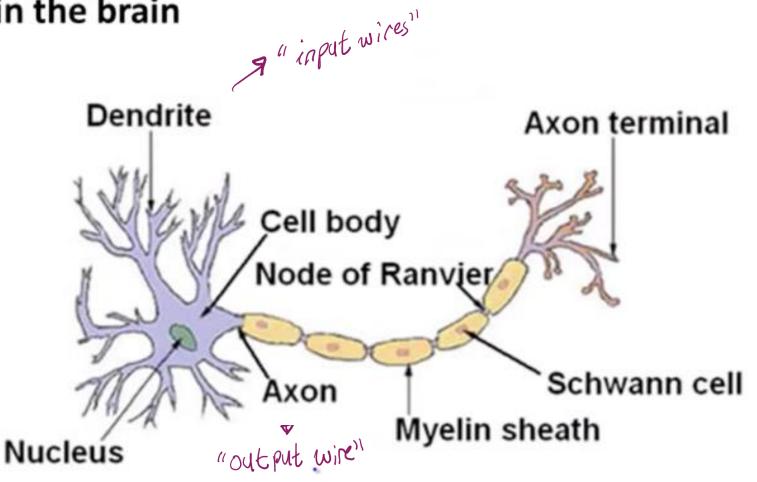
Human echolocation (sonar)



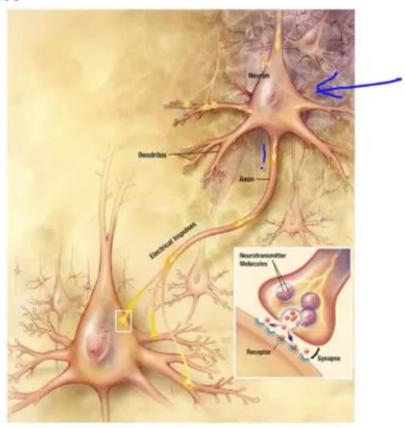
Implanting a 3rd eye

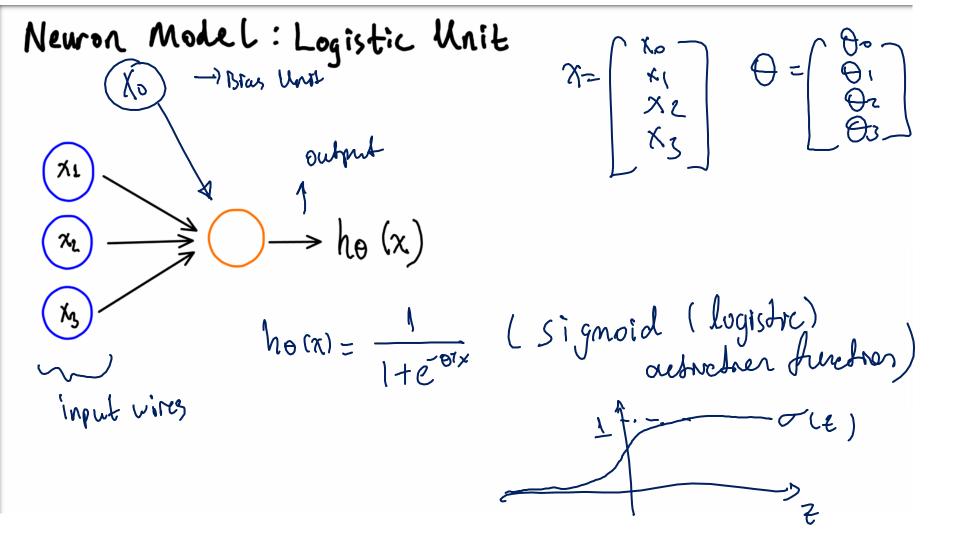
Neural Networks Model Representation I

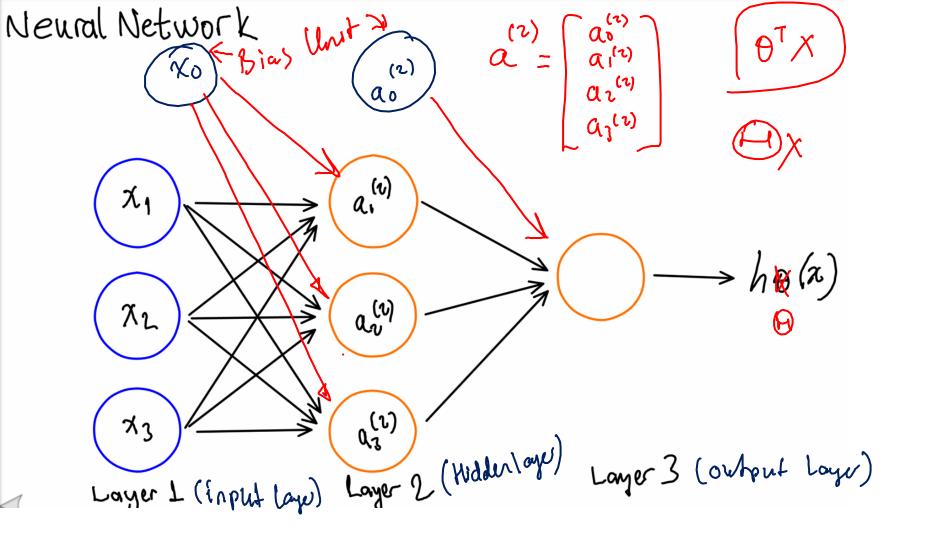
Neuron in the brain



Neurons in the brain







 $\Rightarrow a_3^{(2)} = g(\Theta_{30}^{(1)}x_0 + \Theta_{31}^{(1)}x_1 + \Theta_{32}^{(1)}x_2 + \Theta_{33}^{(1)}x_3)$ $h_{\Theta}(x) = a_1^{(3)} = g(\Theta_{10}^{(2)} a_0^{(2)} + \Theta_{11}^{(2)} a_1^{(2)} + \Theta_{12}^{(2)} a_2^{(2)}$

If network has s_j units in layer j, s_{j+1} units in layer j+1, then $\Theta^{(j)}$

will be of dimension $s_{j+1} \times (s_j + 1)$.

Consider the following neural network: Layer 1 s₁=2 Layer 2 $s_2 = 4$

What is the dimension of $\Theta^{(1)}$ (Hint: add a bias unit to the input and hidden layers)?

$$Sjt1 \times USjt1)$$

$$Sv \times USjt1)$$

$$Sv \times USjt1)$$

$$U \times USjt1$$

$$U \times USjt1$$

$$U \times USjt1$$

$$U \times USjt1$$

Model Representation I

Let's examine how we will represent a hypothesis function using neural networks. At a very simple level, neurons are basically computational units that take inputs (**dendrites**) as electrical inputs (called "spikes") that are channeled to outputs (**axons**). In our model, our dendrites are like the input features $x_1 \cdots x_n$, and the output is the result of our hypothesis function. In this model our x_0 input node is sometimes called the "bias unit." It is always equal to 1. In neural networks, we use the same logistic function as in classification, $\frac{1}{1+e^{-\theta^2x}}$, yet we sometimes call it a sigmoid (logistic) **activation** function. In this situation, our "theta" parameters are sometimes called "weights".

Visually, a simplistic representation looks like:

$$\begin{bmatrix} x_0 \\ x_1 \\ x_2 \end{bmatrix} \to [] \to h_{\theta}(x)$$

Our input nodes (layer 1), also known as the "input layer", go into another node (layer 2), which finally outputs the hypothesis function, known as the "output layer".

We can have intermediate layers of nodes between the input and output layers called the "hidden layers."

In this example, we label these intermediate or "hidden" layer nodes $a_0^2\cdots a_n^2$ and call them "activation units."

$$a_i^{(j)} = \text{"activation" of unit } i \text{ in layer } j$$

 $\Theta^{(j)}$ = matrix of weights controlling function mapping from layer j to layer j+1

If we had one hidden layer, it would look like:

$$\begin{bmatrix} x_0 \\ x_1 \\ x_2 \\ x_3 \end{bmatrix} \rightarrow \begin{bmatrix} a_1^{(2)} \\ a_2^{(2)} \\ a_3^{(2)} \end{bmatrix} \rightarrow h_{\theta}(x)$$

The values for each of the "activation" nodes is obtained as follows:

$$a_{1}^{(2)} = g(\Theta_{10}^{(1)}x_{0} + \Theta_{11}^{(1)}x_{1} + \Theta_{12}^{(1)}x_{2} + \Theta_{13}^{(1)}x_{3})$$

$$a_{2}^{(2)} = g(\Theta_{20}^{(1)}x_{0} + \Theta_{21}^{(1)}x_{1} + \Theta_{22}^{(1)}x_{2} + \Theta_{23}^{(1)}x_{3})$$

$$a_{3}^{(2)} = g(\Theta_{30}^{(1)}x_{0} + \Theta_{31}^{(1)}x_{1} + \Theta_{32}^{(1)}x_{2} + \Theta_{33}^{(1)}x_{3})$$

$$h_{\Theta}(x) = a_{1}^{(3)} = g(\Theta_{10}^{(2)}a_{0}^{(2)} + \Theta_{11}^{(2)}a_{1}^{(2)} + \Theta_{12}^{(2)}a_{2}^{(2)} + \Theta_{13}^{(2)}a_{3}^{(2)})$$

to our inputs to obtain the value for one activation node. Our hypothesis output is the logistic function applied to the sum of the values of our activation nodes, which have been multiplied by yet another parameter matrix $\Theta^{(2)}$ containing the weights for our second layer of nodes.

This is saying that we compute our activation nodes by using a 3×4 matrix of parameters. We apply each row of the parameters

The dimensions of these matrices of weights is determined as follows:

Each layer gets its own matrix of weights, $\Theta^{(j)}$.

If network has s_i units in layer j and s_{i+1} units in layer j+1, then $\Theta^{(j)}$ will be of dimension $s_{i+1} \times (s_i+1)$.

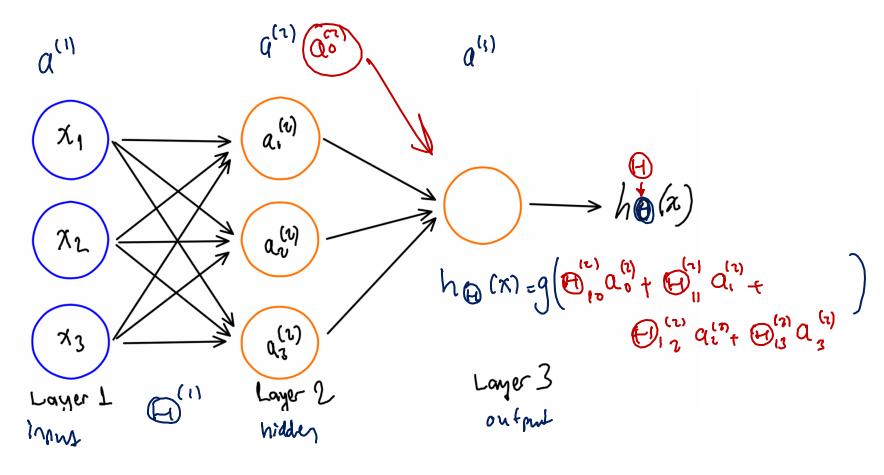
The +1 comes from the addition in $\Theta^{(j)}$ of the "bias nodes," x_0 and $\Theta^{(j)}_0$. In other words the output nodes will not include the

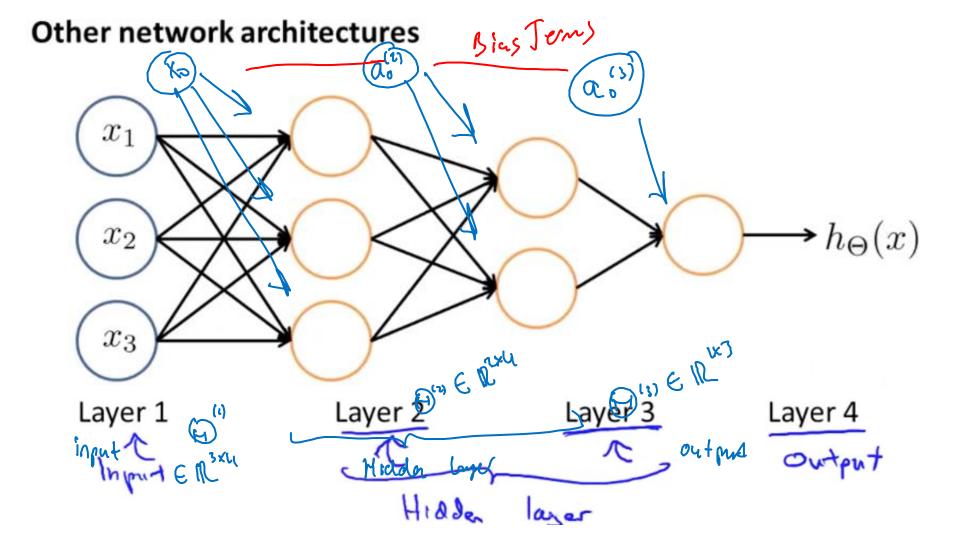
bias nodes while the inputs will. The following image summarizes our model representation:

Neural Networks Model Representation II

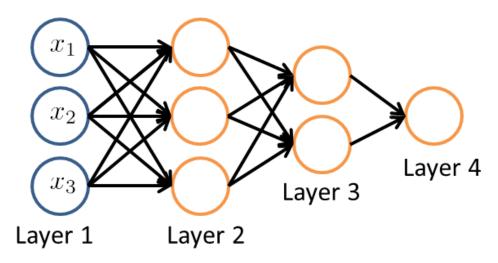
Forward Propagation: Vectorized Implementation
$$\chi = \begin{pmatrix} \chi_1 \\ \chi_2 \\ \chi_3 \end{pmatrix} = \begin{pmatrix} \chi_1 \\ \chi_3 \\ \chi_3 \end{pmatrix} = \begin{pmatrix} \chi_1 \\ \chi_2 \\ \chi_3 \end{pmatrix} = \begin{pmatrix} \chi_1 \\ \chi_3 \\ \chi_3 \end{pmatrix} = \begin{pmatrix} \chi_1 \\ \chi_1 \\ \chi_3 \end{pmatrix} = \begin{pmatrix} \chi_1 \\ \chi_1 \\ \chi_2 \end{pmatrix} = \begin{pmatrix} \chi_1 \\ \chi_1 \\ \chi_3 \end{pmatrix} = \begin{pmatrix} \chi_1 \\ \chi_1 \\ \chi_2 \end{pmatrix} = \begin{pmatrix} \chi_1 \\ \chi_1 \\ \chi_1 \\ \chi_2 \end{pmatrix} = \begin{pmatrix} \chi_1 \\ \chi_1 \\ \chi_1 \\ \chi_2 \end{pmatrix} = \begin{pmatrix} \chi_1 \\ \chi_1 \\ \chi_1 \\ \chi_2 \end{pmatrix} = \begin{pmatrix} \chi_1 \\ \chi_1 \\ \chi_1 \\ \chi_1 \\ \chi_2 \end{pmatrix} = \begin{pmatrix} \chi_1 \\ \chi_1 \\ \chi_1 \\ \chi_1 \\$$

Neural Network learning its own features





Consider the network:



Let $a^{(1)}=x\in\mathbb{R}^{n+1}$ denote the input (with $a_0^{(1)}=1$).

How would you compute $a^{(2)}$?

$$\bigcirc a^{(2)} = \Theta^{(1)}a^{(1)}$$

$$\bigcirc z^{(2)} = \Theta^{(2)}a^{(1)}; \ a^{(2)} = g(z^{(2)})$$

$$\bigcirc \ z^{(2)} = \Theta^{(2)}g(a^{(1)}); \ a^{(2)} = g(z^{(2)})$$

Model Representation II

To re-iterate, the following is an example of a neural network:

$$a_{1}^{(2)} = g(\Theta_{10}^{(1)}x_{0} + \Theta_{11}^{(1)}x_{1} + \Theta_{12}^{(1)}x_{2} + \Theta_{13}^{(1)}x_{3})$$

$$a_{2}^{(2)} = g(\Theta_{20}^{(1)}x_{0} + \Theta_{21}^{(1)}x_{1} + \Theta_{22}^{(1)}x_{2} + \Theta_{23}^{(1)}x_{3})$$

$$a_{3}^{(2)} = g(\Theta_{30}^{(1)}x_{0} + \Theta_{31}^{(1)}x_{1} + \Theta_{32}^{(1)}x_{2} + \Theta_{33}^{(1)}x_{3})$$

$$h_{\Theta}(x) = a_{1}^{(3)} = g(\Theta_{10}^{(2)}a_{0}^{(2)} + \Theta_{11}^{(2)}a_{1}^{(2)} + \Theta_{12}^{(2)}a_{2}^{(2)} + \Theta_{13}^{(2)}a_{3}^{(2)})$$

In this section we'll do a vectorized implementation of the above functions. We're going to define a new variable $z_k^{(j)}$ that encompasses the parameters inside our g function. In our previous example if we replaced by the variable z for all the parameters we would get:

$$a_1^{(2)} = g(z_1^{(2)})$$

 $a_2^{(2)} = g(z_2^{(2)})$
 $a_3^{(2)} = g(z_3^{(2)})$

The vector representation of x and z^j is: $\begin{bmatrix} z_1^{(j)} \end{bmatrix}$

$$x = \begin{bmatrix} x_0 \\ x_1 \\ \dots \\ x_n \end{bmatrix} z^{(j)} = \begin{bmatrix} z_1^{(j)} \\ z_2^{(j)} \\ \dots \\ z_n^{(j)} \end{bmatrix}$$

Setting $x=a^{(1)}$, we can rewrite the equation as:

$$z^{(j)} = \Theta^{(j-1)}a^{(j-1)}$$
We are multiplying our matrix $\Theta^{(j-1)}$ with dimensions $s_i \times (n+1)$ (where s_i is the number of our activation nodes) by our

We are multiplying our matrix $\Theta^{(j-1)}$ with dimensions $s_j \times (n+1)$ (where s_j is the number of our activation nodes) by our vector $a^{(j-1)}$ with height (n+1). This gives us our vector $z^{(j)}$ with height s_j . Now we can get a vector of our activation nodes for layer j as follows:

$$a^{(j)}=g(z^{(j)})$$

Where our function g can be applied element-wise to our vector $z^{(j)}$.

We can then add a bias unit (equal to 1) to layer j after we have computed $a^{(j)}$. This will be element $a_0^{(j)}$ and will be equal to 1. To compute our final hypothesis, let's first compute another z vector:

$$z^{(j+1)} = \Theta^{(j)}a^{(j)}$$

We are multiplying our matrix $\Theta^{(j-1)}$ with dimensions $s_j \times (n+1)$ (where s_j is the number of our activation nodes) by our vector $a^{(j-1)}$ with height (n+1). This gives us our vector $z^{(j)}$ with height s_j . Now we can get a vector of our activation nodes for layer j as follows:

$$a^{(j)} = g(z^{(j)})$$

Where our function g can be applied element-wise to our vector $z^{(j)}$.

We can then add a bias unit (equal to 1) to layer j after we have computed $a^{(j)}$. This will be element $a_0^{(j)}$ and will be equal to 1. To compute our final hypothesis, let's first compute another z vector:

$$z^{(j+1)} = \Theta^{(j)}a^{(j)}$$

We get this final z vector by multiplying the next theta matrix after $\Theta^{(j-1)}$ with the values of all the activation nodes we just got. This last theta matrix $\Theta^{(j)}$ will have only **one row** which is multiplied by one column $a^{(j)}$ so that our result is a single number. We then get our final result with:

$$h_{\Theta}(x) = a^{(j+1)} = g(z^{(j+1)})$$

Notice that in this **last step**, between layer j and layer j+1, we are doing **exactly the same thing** as we did in logistic regression. Adding all these intermediate layers in neural networks allows us to more elegantly produce interesting and more complex non-linear hypotheses.