# **Insurance analytics**

### Neural networks

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## Acknowledgement

- ▶ Some of the figures in this presentation are taken from
  - Michael A. Nielsen (2015). Neural networks and deep learning. Determination Press.
  - all the beautiful work by prof. Taylor Arnold, in particular Chapter 8 of the book *A Computational Approach to Statistical Learning*.

# Today's mission

- ► Today's mission:
  - de-mystify neural networks
  - sketch different types of neural networks and their applications
  - a discussion of specific considerations to keep in mind when using these predictive modeling techniques with frequency/severity data.

# A famous challenge

Recognizing handwritten digits

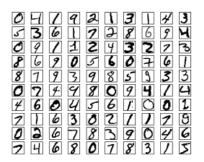
Consider this

# 504192

- ▶ Humans have a primary visual cortex  $(V_1)$  with 140M neurons, with 10s billions connections.
- ▶ Next to  $V_1$ , also  $V_2$ ,  $V_3$ ,  $V_4$  and  $V_5$ , doing complex image processing.
- A supercomputer in our head!

## A famous challenge

Recognizing handwritten digits



The goal: infer rules for recognizing handwritten digits.

The famous MNIST (Modified National Institute of Standards and Technology) database with handwritten digits commonly used for training various image processing systems.

## A famous challenge

Recognizing handwritten digits



The winner: well-designed neural nets classify 9 979 out of 10 000 images correctly (in 2013).

## Types of neural networks

What's in a name?

- ▶ Different types of neural networks and their applications:
  - ANN: Artificial Neural Network

for regression and classification problems, with vectors as input data

- CNN: Convolutional Neural Network
  - for image processing, image/face/ $\dots$  recognition, with images as input data
- RNN: Recurrent Neural Network

for sequential data such as text or time series.

## Types of neural networks

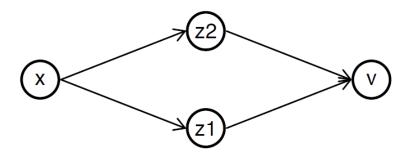
#### A bit of history

- ▶ 1943: McCulloch and Pitts (1943) with a computational model for neural networks based on maths and algorithms. Paved the way for neural network research in artificial intelligence.
- ▶ 1958: perceptron algorithm for pattern recognition by Rosenblatt.
- ▶ 1969: discovery of two fundamental problems with the computational machines that processed neural networks (e.g. computer power) by Minsky and Papert.
- ▶ 1986: backpropagation algorithm published in Nature.
- ▶ 1990 2010: not much happening.
- ▶ 2010 ... : Al boom, driven by (deep) neural networks. Many factors play a role: power of CNNs and RNNs recognized, more computer power, higher demand due to more and more complex data.

A simple neural network

- ► De-mystify artificial neural networks (ANNs):
  - a collection of inter-woven linear models
  - extending linear approaches to detect non-linear interactions in high-dimensional data.

A simple neural network



The goal: predict a scalar response y from scalar input x.

A simple neural network

- ► Some terminology:
  - x is the input layer
  - v is output layer
  - middle layers are hidden layers
  - four neurons: x,  $z_1$ ,  $z_2$  and v.

A simple neural network

First, apply two independent linear models:

$$z_1 = b_1 + x \cdot w_1$$
  
$$z_2 = b_2 + x \cdot w_2,$$

using four parameters: two intercepts and two slopes.

Next, construct another linear model with the  $z_i$  as inputs:

$$\hat{y} := v = b_3 + z_1 \cdot u_1 + z_2 \cdot u_2.$$

A simple neural network

Putting it all together:

$$v = b_3 + z_1 \cdot u_1 + z_2 \cdot u_2$$

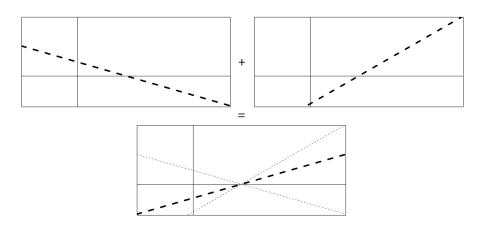
$$= b_3 + (b_1 + x \cdot w_1) \cdot u_1 + (b_2 + x \cdot w_2) \cdot u_2$$

$$= (b_3 + u_1 \cdot b_1 + u_2 \cdot b_2) + (w_1 \cdot u_1 + w_2 \cdot u_2) \cdot x$$

$$= (intercept) + (slope) \cdot x.$$

- Model is over-parametrized, with infinitely many ways to describe same model.
- Essentially, still a linear model!

A simple neural network



#### Activation function

ightharpoonup Capture non-linear relationships between x and v by replacing

$$v = b_3 + z_1 \cdot u_1 + z_2 \cdot u_2.$$

with

$$v = b_3 + \sigma(z_1) \cdot u_1 + \sigma(z_2) \cdot u_2 = b_3 + \sigma(b_1 + x \cdot w_1) \cdot u_1 + \sigma(b_2 + x \cdot w_2) \cdot u_2,$$

where  $\sigma(.)$  is an activation function, a mapping from  $\mathbb{R}$  to  $\mathbb{R}$ .

#### Activation function

ightharpoonup Capture non-linear relationships between x and v:

$$v = b_3 + \sigma(z_1) \cdot u_1 + \sigma(z_2) \cdot u_2,$$

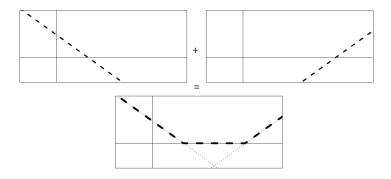
where  $\sigma(.)$  is an activation function, a mapping from  $\mathbb{R}$  to  $\mathbb{R}$ .

For example, the rectified linear unit (ReLU) activation function:

$$ReLU(x) = \begin{cases} x, & \text{if } x \ge 0 \\ 0, & \text{otherwise.} \end{cases}$$

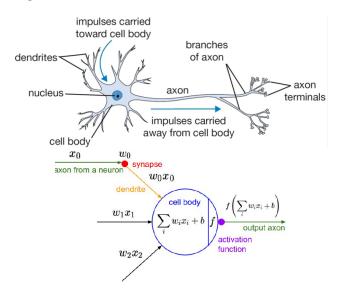
Adding an activation function greatly increases the set of possible relations between x and v!

#### Activation function



More activation functions: sigmoid, hyperbolic tan, leaky rectified linear unit, maxout.

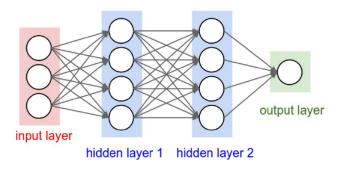
#### Artificial vs biological



#### The architecture

- Artificial Neural networks (NNs):
  - · a collection of neurons
  - organized into an ordered set of layers
  - directed connections pass signals between neurons in adjacent layers
  - to train: update parameters describing the connections by minimizing loss function over training data
  - to predict: pass  $x_i$  to first layer, output of final layer is  $\hat{y}_i$ .
- ► The network is *dense* if each neuron in a layer receives an input from all the neurons present in the previous layer; *densely connected*.

The architecture



A basic neural network. Source: http://blog.christianperone.com

This is a feedforward neural network - no loops!

- Using the NN language:
  - intercept called the bias
  - slopes called weights
  - L layers in total, with input layer denoted as layer 0
  - use a (from activation) to denote the output of a given layer.
- ▶ Let's start with a single layer network (called an artificial neuron).

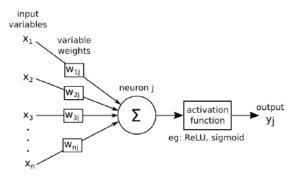
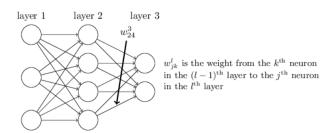
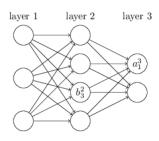


Figure 1: source: andrewjames turner.co.uk

- Use  $w_{ik}^I$  to denote the weight for the connection:
  - from neuron k in layer (l-1)
  - to neuron j in layer l.



- ▶ Use
  - $b_j^I$  for the bias of neuron j in layer I
  - $a_i^I$  for the activation of neuron j in layer I.



#### **Terminology**

• *d*-dimensional inputs

$$a^0 = x$$

• weight  $w_{i,k}^1$  on kth neuron in layer 0 within jth neuron in layer 1

$$w_j^1 = (w_{j,1}^1, \dots, w_{j,d}^1)$$

• with bias term  $b_i^1$ 

$$z_j^1 = w_{j,1}^1 \cdot x_1 + \ldots + w_{j,d}^1 \cdot x_d + b_j^1$$
  
=  $\langle w_j^1, x \rangle + b_j^1$ ,

apply activation function

$$a_i^1 = \sigma(z_i^1),$$

the output of neuron j in layer 1.

#### Terminology

▶ With *L* layers in total: (in vectorized form)

$$z' = \langle w', a'^{-1} \rangle + b'$$
 (the weighted input)  $a' = \sigma(z')$ ,

for I from 1 up to (and including) L.

► Finally,

$$a^L = \hat{y} = \tilde{\sigma}(z^L),$$

with examples of output activations: identity (regression), sigmoid (binary classification), softmax (muti-class output).

Basic idea - GD

► We want to find

$$\min_{w} f(w)$$
,

then gradient descent updates as follows (cfr. lecture on *boosting methods*)

$$w_{\text{new}} = w_{\text{old}} - \eta \cdot \nabla_w f(w_{\text{old}}),$$

with learning rate  $\eta$ . Move in the direction the function locally decreases the fastest!

► A good choice for NNs because faster second-order methods involve the Hessian and this is infeasible when having tons of parameters.

#### From GD to SGD

▶ Let  $\mathcal{L}(w; y_i, x_i)$  be the loss function and w the parameters. For example,

$$\mathcal{L}(w; y_i, x_i) = \frac{1}{2n} \sum_i (\hat{y}_i(w) - y_i)^2$$
$$= \frac{1}{n} \sum_i \mathcal{L}_i(w; y_i, x_i).$$

▶ With the gradient descent update rule

$$w_{\text{new}} = w_{\text{old}} - \eta \cdot \nabla_w \mathcal{L}(w),$$

where 
$$\nabla_w \mathcal{L}(w) = \frac{1}{n} \sum_i \nabla_w \mathcal{L}_i$$
.

#### Mini-batches

Stochastic gradient descent picks randomly m training inputs  $x_1, \ldots, x_m$ , a mini-batch:

$$\frac{\sum_{j=1}^{m} \nabla_{w} \mathcal{L}_{j}}{m} \approx \frac{\sum_{i} \nabla_{w} \mathcal{L}_{i}}{n} = \nabla_{w} \mathcal{L}.$$

► Thus,

$$\nabla_{w}\mathcal{L} \approx \frac{1}{m}\sum_{i=1}^{m}\nabla\mathcal{L}_{j}(w).$$

Estimate the gradient only on a small subset of the entire training set.

- ▶ Partition input randomly into disjoint groups  $M_1, M_2, \ldots, M_{n/m}$ .
- Updates:

$$w_{k+1} = w_k - \frac{\eta}{m} \sum_{i \in M_1} \nabla \mathcal{L}_i$$

$$w_{k+2} = w_{k+1} - \frac{\eta}{m} \sum_{i \in M_2} \nabla \mathcal{L}_i$$

$$\vdots$$

$$w_{k+n/m+1} = w_{k+n/m} - \frac{\eta}{m} \sum_{i \in M_{n/m}} \nabla \mathcal{L}_i.$$

Speed up the process of doing gradient descent. Going through the entire data set is called an epoch.

- ► Compute the gradient of the loss function wrt all trainable parameters:
  - tons of parameters
  - need for efficient algorithm to calculate gradient
  - generic algorithm usable for arbitrary number of layers and neurons in each layer.
- ▶ The strategy (Rumelhart et al., 1986, Nature)
  - backwards propagation of errors or backpropagation
  - uses chain rule for derivatives.

 $\blacktriangleright$  With a loss function  $\mathcal{L}$ , e.g. squared error loss

$$\mathcal{L}(y,a^L) = \frac{1}{2n}(y-a^L)^2.$$

▶ Equations defining backpropagation: (recall:  $z_j^l = \sum_k w_{jk}^l a_k^{l-1} + b_j^l$ ) starting point - gradient terms of  $b_i^l$ 

$$\frac{\partial \mathcal{L}}{\partial b_{j}^{l}} = \frac{\partial \mathcal{L}}{\partial z_{j}^{l}} \cdot \frac{\partial z_{j}^{l}}{\partial b_{j}^{l}} 
= \frac{\partial \mathcal{L}}{\partial z_{i}^{l}} \cdot 1 = \frac{\partial \mathcal{L}}{\partial z_{i}^{l}}.$$

Thus, problem centers around derivatives wrt  $z^{l}$ !

 $\blacktriangleright$  With a loss function  $\mathcal{L}$ , e.g. squared error loss

$$\mathcal{L}(y, a^L) = \frac{1}{2n}(y - a^L)^2.$$

▶ Equations defining backpropagation: (recall:  $z_j^l = \sum_k w_{jk}^l a_k^{l-1} + b_j^l$ )

starting point - gradient terms of weights  $w_{jk}^I$ 

$$\frac{\partial \mathcal{L}}{\partial w_{jk}^{l}} = \frac{\partial \mathcal{L}}{\partial z_{j}^{l}} \cdot \frac{\partial z_{j}^{l}}{\partial w_{jk}^{l}}$$
$$= \frac{\partial \mathcal{L}}{\partial z_{i}^{l}} \cdot a_{k}^{l-1}.$$

Thus, problem centers around derivatives wrt  $z^{l}$ !

The derivatives in layer L are straightforward to calculate (with  $a_i^L = \sigma(z_i^L)$ )

$$\frac{\partial \mathcal{L}}{\partial z_{j}^{L}} = \sum_{k} \frac{\partial \mathcal{L}}{\partial a_{k}^{L}} \cdot \frac{\partial a_{k}^{L}}{\partial z_{j}^{L}} 
= \frac{\partial \mathcal{L}}{\partial a_{j}^{L}} \cdot \frac{\partial a_{j}^{L}}{\partial z_{j}^{L}} 
= \frac{\partial \mathcal{L}}{\partial a_{k}^{L}} \cdot \sigma'(z_{j}^{L}).$$

An equation for the error in the output layer!

- ► The derivatives wrt z<sup>l</sup> can be written as a function of derivatives wrt z<sup>l+1</sup>.
- ► Using the chain rule

$$\frac{\partial \mathcal{L}}{\partial z_{j}^{l}} = \sum_{k} \frac{\partial \mathcal{L}}{\partial z_{k}^{l+1}} \cdot \frac{\partial z_{k}^{l+1}}{\partial z_{j}^{l}} 
= \sum_{k} \frac{\partial \mathcal{L}}{\partial z_{k}^{l+1}} \cdot w_{kj}^{l+1} \sigma'(z_{j}^{l}),$$

where we use

$$z_k^{l+1} = \sum_i w_{kj}^{l+1} a_j^l + b_k^{l+1} = \sum_i w_{kj}^{l+1} \sigma(z_j^l) + b_k^{l+1}.$$

- ▶ The vanilla implementation of the backpropagation alogrithm:
  - 1. Input x, set  $a^0$  for the input layer.
  - 2. Feedforward: for each l = 2, 3, ..., L compute

$$z' = \langle w', a'^{-1} \rangle + b' \text{ and } a' = \sigma(z').$$

- 3. Output error: compute  $\frac{\partial \mathcal{L}}{\partial z_i^L} = \frac{\partial \mathcal{L}}{\partial a_i^L} \cdot \sigma'(z_j^L)$ .
- 4. Backpropagate the error: for each  $l = L 1, \ldots, 2$  compute  $\frac{\partial \mathcal{L}}{\partial z_i^l} = \sum_k \frac{\partial \mathcal{L}}{\partial z_k^{l+1}} \cdot w_{kj}^{l+1} \sigma'(z_j^l)$ .
- 5. Output: the gradient of the loss function  $\frac{\partial \mathcal{L}}{\partial w_{ik}^{l}} = \frac{\partial \mathcal{L}}{\partial z_{i}^{l}} \cdot a_{k}^{l-1}$  and  $\frac{\partial \mathcal{L}}{\partial b_{i}^{l}} = \frac{\partial \mathcal{L}}{\partial z_{i}^{l}}$ .

- Several improvements covered in the literature
  - to reduce overfitting
  - to reduce getting stuck in local saddle points
  - to converge in a smaller number of epochs.

Reduce overfitting

- ► Methods to reduce overfitting:
  - different weight initialization
  - early stopping
  - regularization
  - dropout.

Reduce overfitting: weight initialization, early stopping

- ▶ Update the weight initialization ~> might drastically improve performance of a model with many layers.
- ► Early stopping to prevent overfitting:
  - calculate validation performance after each epoch
  - stop when this no longer improves
  - NNs are motivated by an optimization problem, but do not attempt to solve the optimization task.

Reduce overfitting: regularization

- Prevent overfitting via regularization (cfr. lecture on Lasso and friends)
  - add (e.g.)  $\ell_2$ -norm

$$f_{\lambda}(w,b) = f(w,b) + \frac{\lambda}{2} \cdot ||w||_2^2,$$

with weights w and bias terms b, where only weights are penalized

with gradient

$$\nabla_w f_{\lambda} = \nabla_w f(w, b) + \lambda \cdot w$$

such that

$$\begin{array}{lll} w_{\mathsf{new}} & \leftarrow & w_{\mathsf{old}} - \eta \cdot \nabla_{w} f_{\lambda}(w_{\mathsf{old}}) \\ & \leftarrow & w_{\mathsf{old}} - \eta \cdot [\nabla_{w} f(w_{\mathsf{old}}) + \lambda \cdot w] \\ & \leftarrow & [1 - \eta \cdot \lambda] \cdot w_{\mathsf{old}} - \eta \cdot f(w_{\mathsf{old}}). \end{array}$$

Reduce overfitting: dropout

### ► Dropout:

- randomly set activations to zero, with fixed probability p
- both in forward propagation as well as backpropagation
- only in training, all nodes turned on during prediction.

# Improving SGD and regularization Dropout

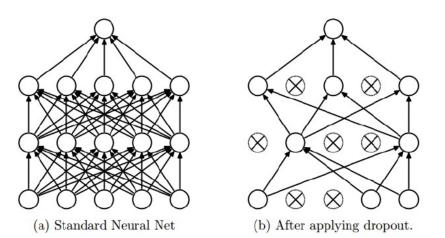


Figure 5: Dropout - source: http://blog.christianperone.com/

Saddle points & faster convergence

- ▶ There is a fast proliferation of improved optimization algorithms.
- Examples:
  - momentum
  - RMSprop
  - · adagrad, adamdelta
  - adam, adamax
  - natural gradient descent

# Where to finetune your ANN?

- ► A list of tuning parameters/architecture choices:
  - learning rate  $\eta$
  - regularization parameter  $\lambda$
  - number of epochs
  - · mini-batch size
  - number of layers
  - number of hidden neurons per layer
  - activation functions, choice of optimization strategy.

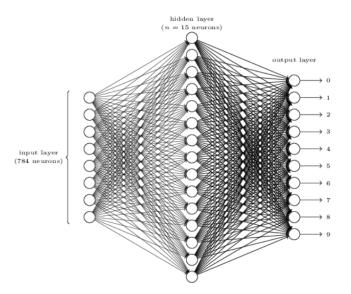
### Classification with Artificial neural networks

Back to the famous challenge

- ▶ The goal: try to recognize individual hand-written digits
  - take e.g. 28 by 28 greyscale image
  - $784 = 28 \times 28$  input neurons, with intensities between 0 (white) and 1 (black)
  - output layer has 10 layers
    - neuron with highest activation value fires.

### Classification with artificial neural networks

#### Back to the famous challenge



### Classification with artificial neural networks

- Use a multi-valued output layer for classification tasks.
- One-hot encoding applied to output vector y

$$\begin{pmatrix} 2\\4\\\vdots\\1 \end{pmatrix} \rightarrow \begin{pmatrix} 0 & 1 & 0 & 0 & 0\\0 & 0 & 0 & 1 & 0\\\vdots & \vdots & \vdots & \vdots\\1 & 0 & 0 & 0 & 0 \end{pmatrix}.$$

Use softmax function as activation in final layer

$$a_j^L = \operatorname{softmax}(z_j^L)$$
  
=  $\frac{e^{z_j}}{\sum_k e^{z_k}}$ ,

for the *j*th output neuron. The last layer is then easily interpreted as sequence of probabilities.

### Classification with artificial neural networks

Instead of using squared error loss, use categorical cross-entropy

$$\mathcal{L}(a^L, y) = -\frac{1}{n} \sum_{i} \sum_{k} y_{ik} \cdot \log(a_k^L),$$

as the loss function, where i runs over training data and k runs over the class labels.

- ▶ Think: softmax output layer and negative log-likelihood cost function.
- ► Work out the derivatives to set-up the backpropagation with this loss function.

The motivation

- Prediction tasks with images as inputs are a popular application of NNs.
- With a limited number of input pixels, put a weight on each pixel and use ANN.
- ▶ With large images use convolutional layers.

#### A convolution?

 $\triangleright$  The discrete convolution between f and g is

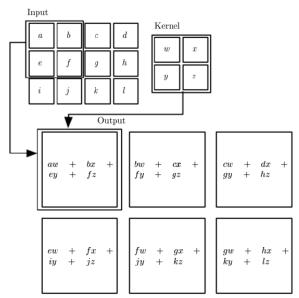
$$(f*g)(x) = \sum_{t} f(t) \cdot g(x+t).$$

▶ With 2-dimensional signals (e.g. images), consider 2D-convolutions

$$(K*I)(i,j) = \sum_{m,n} K(m,n) \cdot I(i+m,j+n),$$

with K a convolutional kernel applied to a 2D signal (or image) I.

#### A convolution pictured



- ▶ With large images use convolutional layers:
  - · apply small set of weights to subsections of the image
  - same weights across the image
  - use a kernel matrix, e.g. for a black and white image

$$K = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}.$$

Take pixel value and subtract from this the pixel value to its immediate lower right.

- ► In a CNN:
  - apply several such kernels in a layer
  - with weights learned during training of the algorithm
  - different convolutions pick up different features (e.g. edges, texture, basic object types).
- Pooling layer: simplify information in output from convolutional layer.
- ▶ With input image in color: kernel matrix *K* is 3-dimensional array with weights applied to each color channel.

### Recurrent neural networks

- ▶ To infer sequential data such as text or time series.
- ► Mathematically,

$$h_t = Wx_t + b + Uh_{t-1}$$
  
=  $Wx_t + b + UWx_{t-1} + Ub + U^2h_{t-2}$   
= ...

### Recurrent neural networks

#### Unrolled

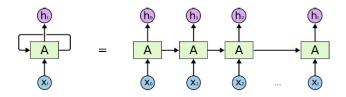


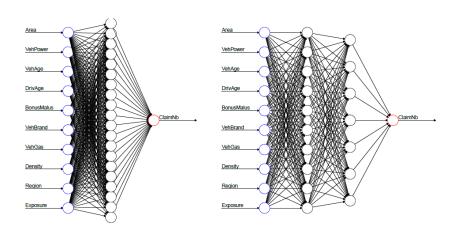
Figure 19: Unrolled representation of a RNN. Source : Understanding LSTM Networks by Christopher Olah - http://colah.github.io/posts/2015-08-Understanding-LSTMs/

► Starting from the Poisson regression model

$$N_i \sim POI(\lambda(x_i, d_i)),$$

with  $d_i$  exposure and  $x_i$  the features for insured i.

- ► Build e.g.
  - a fully-connected single hidden layer feed-forward neural network with 20 hidden neurons for modelling the regression function
  - a more general architecture with K hidden layers.



A shallow (left) and a deep (right) neural network.

- Some comments:
  - output layer has a single neuron, and exponential activation function
  - · some preprocessing steps:
    - ordered categorical features as continuous
    - dummy coding or one-hot encoding to construct binary representation of nominal features
    - put continuous covariates on the same scale
  - use Poisson loss function in keras.

#### Listing 3: R script for fitting networks in Keras

```
library(keras)
2
   model <- keras_model_sequential()</pre>
    model %>%
5
      layer_dense(units = q1, activation = 'tanh', input_shape = c(ncol(Xlearn))) %>%
      layer dense(units = 1, activation = k exp)
6
7
8
    summary (model)
g
10
   model %>% compile(
11
      loss = 'poisson'.
12
     optimizer ='sgd'
13
   )
14
15
   fit <- model %>% fit(Xlearn, learn$ClaimNb, epochs=100, batch_size=10000)
```

keras implementation of a neural network for frequencies.

K. Antonio, KU Leuven & UvA Yes we CANN! 59/60

- Explore the following links
  - Case study: French MTPL claims
  - Insights from inside neural networks
  - Data analytics for non-life insurance pricing
  - Editorial: Yes we CANN!

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