Building LLM for Statisticians

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LLM for Statisticians

- For statisticians like myself, LLMs (Large Language Models) are fairly easy to understand in principle as they are just machine learning models with a few tricks:
 - Nonlinear activation functions allow complex non-linear models with an overall mostly linear (additive) structure
 - Thus easily differentiable which allows straightforward optimization via gradient descent
 - Layered structure so parameters can be updated in manageable blocks (back-propagation)
 - Large matrix (tensor) operations are amenable to parallel execution and perfect for GPUs
- Still it is a bit of a challenge for a statistician like me to understand the implementation of LLMs due to esoteric terminology (what the heck is "attention"?), Python codes, and math (either too much or too little, and in unfamiliar notations).
- Build a LLM from Scratch is an excellent book that allows one to take a peek at how LLM is actually implemented. Still it uses Python and very little math. Here my notes use R instead with some math, that's easier for myself to understand.

From GPT to ChatGPT

Large Language Model (LLM) is a (large) neural network (NN) model for next-word prediction.

- 1. The input is sequence of tokens (words, punctuation, and possibly other special context token like end-of-text) of length *context size* (typically 256).
 - tokenization: breaking down input texts into individual tokens
 - encoding: converting tokens into integers (token IDs) by looking up inthe dictionary
 - embedding: converting one token ID to a vector of real numbers
- 2. Pre-training takes a large corpus of text to create a base or /foundation/ model, such as GPT (Generative Pretrained Transformers). This is called self-supervised training as the training data (the target is the next word/token) is created on-the-fly by a data loader. This model is capable of text completion (zero-shot) and has limited few-shot capabilities, meaning learning to perform new tasks based on only a few examples.
- 3. Fine-tuning involves further training the LLM on labeled data. Two popular categories are /instruction/ (instructions and answers) and /classification/ (texts and labels) fine-tuning. ChatGPT is such a fine-tuned model.

LLM in a Nutshell

■ Like all models, LLM is a function that maps the input X (a sequence of words or tokens) to an output Y (a vector of multinomial probability for any token in the "vocabulary").

$$Y = f_{LLM}(X)$$

- In LLM, calculations are done in a step-wise fashion, where each step (or "layer") can be one of the following, and the output of one step is the input of the next.
 - Linear transformation (in matrix form): $f(x) = xA^{T} + b$
 - Ancillary actions like normalization or dropout to improve computational efficiency and numerical stability.
 - Activation functions to introduce non-linearity.
- At each step, the output is a matrix or a tensor (multidimensional matrices, which can also be viewed as a larger block-sparse matrix) where each element is often fancifully referred to as a "neuron" or a "node" in the neural network.
- Despite the simple structure, this model can approximate any function when sufficiently large. The model parameters (to be optimized/trained) are only involved in the linear operations. Thus model fitting is feasible even for very large models.

What's the Trick?

- **Transformer** refers to the two steps of the process, an *encoder* that prepares the input, and an *decoder* that does the prediction. Both encoder and decoder can be large neural networks that are optimized with training. For LLMs the encoder is relatively simple and GPT focuses on the decoder only.
- LLMs like GPT taken into account contextual information, called *attention mechanism*. It's essentially a weighted correlation matrix of the tokens, where the weights are parameters to be optimized via training.
- GPT-3's vocabulary includes V = 50257 tokens, with a maximum context size of 2,049 tokens. It uses an embedding size of d = 12288 dimensions and L = 96 layers. The total number of parameters is:

$$V \times d + L \times (4d^2 + 8d^2 + 4d) + d \times V \approx 2 \times 50257 \times 12288 + 12 \times 96 \times 12288^2 \approx 175 \times 10^9$$

 As the model scales up (increased number of nodes or parameters) over certain threshold, unexpected capabilities arise, such as chatGPT doing math. This is called emergent behavior.

PyTorch and R

- PyTorch has surpassed TensorFlow as the ML library of choice (in 2025).
- torch is a native-R interface for GPU accelerated array computation, general neural network layers and data loaders.

```
install.packages(c("torch", "libtorch", "bench"))
```

 On Apple Silicon, Metal Performance Shaders (MPS) backend provides GPU acceleration for matrix calculation. CUDA is used with NVIDIA GPUs. Here the tensor (an array) is stored in GPU.

```
torch::backends_mps_is_available()
## [1] TRUE
a.mps <- torch::torch_rand(c(1000, 1000), device = "mps")
b.mps <- torch::torch_rand(c(1000, 1000), device = "mps")
a.mps[1:2,1:8]
## torch_tensor
## 0.2847 0.3973 0.9793 0.3278 0.1432 0.4976 0.3536 0.1886
## 0.3032 0.2595 0.7262 0.2954 0.6430 0.2013 0.5535 0.7115
## [ MPSFloatType[2,8] ]</pre>
```

Note that in R, vector and array index starts from 1 (FORTRAN style), versus 0 (C style) for Python.

GPU Acceleration

Tensor on CPU

```
a.cpu <- torch::torch_rand(c(1000, 1000), device = "cpu")
(b.cpu <- torch::torch_rand(c(1000, 1000), device = "cpu"))[1:2, 1:8]

## torch_tensor
## 0.7336 0.2180 0.2343 0.0688 0.1203 0.2866 0.1235 0.9458

## 0.2443 0.5215 0.9779 0.8567 0.3307 0.7691 0.9304 0.1461

## [ CPUFloatType{2,8} ]
```

■ Plain matrix

```
a <- matrix(runif(n = 1000*1000), ncol = 1000)
b <- matrix(runif(n = 1000*1000), ncol = 1000)</pre>
```

Multiplication

```
rbind(bench::mark(c.mps <- torch::torch mm(a.mps, b.mps)).</pre>
     bench::mark(c.cpu <- torch::torch mm(a.cpu, b.cpu)), bench::mark(c <- a %*% b))[, 1:9]
## # A tibble: 3 x 6
    expression
                                           min median 'itr/sec' mem alloc 'gc/sec'
    <bch:expr>
                                         <bch> <bch:t>
                                                           <dbl> <bch:bvt>
                                                                              <fdh1>
## 1 c.mps <- torch::torch_mm(a.mps, b.~ 31us 52.1us
                                                                   13.89KB
                                                         2235.
## 2 c.cpu <- torch::torch mm(a.cpu, b.~ 931us 1.2ms
                                                          731.
                                                                        ΘB
## 3 c <- a %*% b
                                         322ms 327ms
                                                           3.06
                                                                    7.63MB
                                                                                  0
```

More about Torch Tensor

■ Torch tensors are somewhat mysterious R7 (or S7?) class objects, and pointers to Python objects.

```
aa <- torch::torch_tensor(array(c(0.2745, 0.6584, 0.2775, 0.8573, 0.8993, 0.0390, 0.9268, 0.7388, 0.0772, 0.3565, 0.1479, 0.5331, 0.4066, 0.2318, 0.4545, 0.9737), dim = c(1, 2, 2, 4)))

class(aa)
## [1] "torch tensor" "R7"
```

As pointers to Python objects, they have some "methods" available, but some ordinary R (base or from torch library) functions also work.

```
aa$size()
## [1] 1 2 2 4
dim(aa)
## [1] 1 2 2 4
```

■ Two ways of combining tensors: "stack" increases the number of dimensions while "cat" does not. The first line concatenate two $1 \times 2 \times 4 \times 2$ tensors to create a tensor of $1 \times 2 \times 4 \times 4$. The second line stacks them to create a tensor of $2 \times 1 \times 2 \times 4 \times 2$.

```
torch::torch_cat(c(aa$transpose(3, 4), torch::torch_transpose(aa, 3, 4)), dim = 4L)
torch::torch_stack(c(aa$transpose(3, 4), torch::torch_transpose(aa, 3, 4)), dim = 1L)
```

Tokenization

- Tokenization is the process of breaking down raw text into small pieces (tokens, including words, punctuation, etc) and then convert the character strings to integers through by looking up in the vocabulary.
- rtiktoken provides R interface to OpenAI's BPE tokenizer python library tiktoken

```
fname <- "https://raw.githubusercontent.com/rasbt/LLMs-from-scratch/refs/heads/main/cho2/o1_main-chapter-code/the-verdict.txt"
nchar(raw.text <- readr::read_file(fname))
## [1] 20479
enc.text <- rtiktoken::get_tokens(raw.text, model = "gpt2")
length(enc.text)
## [1] 5145</pre>
```

Encoded values:

```
enc.text[1:10]
## [1] 40 367 2885 1464 1807 3619 402 271 10899 2138
```

Converting back, note that it may not identical to the original text as white spaces etc may have been dropped.

```
rtiktoken::decode_tokens(enc.text[1:10], "gpt2")
## [1] "I HAD always thought Jack Gisburn rather"
```

Data Sampling with a Sliding Window

- Inputs and targets are generated automatically in training a LLM through a sliding window approach. An input vector of context size 4 (max.len) is first selected, the target is then shifted by 1 (for next word prediction).
- The next input is shifted for a stride of size 3 (stride). When the stride size equals context size there is no overlap between inputs thus less correlation.

```
n <- length(enc.text)</pre>
idx <- seq(1, n - max.len, by = stride)
tmp <- cbind(enc.text[(1:(n-max.len))]. enc.text[(2:(n-max.len+1))])</pre>
inputs <- targets <- array(o, dim = c(batch.size, max.len))
for (i in seg(batch.size)) {
    this.idx <- idx[i]:(idx[i]*stride)</pre>
    inputs[i,] <- tmp[this.idx.1]
    targets[i.] <- tmp[this.idx.2]</pre>
(inputs.t <- torch::torch tensor(inputs. dtype = torch::torch long()))
## torch tensor
     40 367 2885 1464
    1464 1807 3619
## [ CPULongType{2.4} ]
(targets.t <- torch::torch_tensor(targets, dtype = torch::torch_long()))</pre>
## torch tensor
    367 2885 1464 1807
    1807 3619 402
                       271
    CPULongType{2.4} 1
```

Embeddings

- Embedding creates a continuous vector representation for each token ID, which is necessary for training the deep neural network through backpropagation algorithm.
- These are essentially weights that correspond to the contribution of each token to the bottom layer of the network and are optimized through training.

```
embedding.ex <- torch::nn_embedding(num_embeddings = 6, # size of the vocabulary</pre>
                                    embedding dim = 3) # dimension or number of neural nodes at this layer
embedding.exsweight
## torch tensor
## -0.7695 1.1331 -1.2419
## 1.1693 2.2465 -0.7827
## -1.6537 1.0635 0.2136
## -1.3453 -0.0734 0.6027
## -1.0212 -0.8909 0.0688
## -0.9263 -1.2714 0.5308
## [ CPUFloatType{6.3} ][ requires grad = TRUE ]
input.ex <- torch::torch tensor(c(1, 2, 4), dtype = torch::torch long()) # needs Long integers as indices
embedding.ex(input.ex)
## torch tensor
## -0.7695 1.1331 -1.2419
## 1.1603 2.2465 -0.7827
## -1.3453 -0.0734 0.6027
## [ CPUFloatType{3.3} ][ grad fn = <EmbeddingBackwardo> ]
```

Embedding with Positional Information

First the create embeddings for the tokens IDs:

GPT uses absolute position embeddings (weights for each position).

Position embeddings and token embeddings add up to the input embeddings.

```
input.embeddings <- token.embeddings * pos.embeddings
input.embeddings[1,1,1:5]

## torch_tensor
## 0.2646
## 1.3290
## -0.5708
## 0.3094
## 1.7389
## [ CPUFloatType{5} ][ grad_fn = <SliceBackwardo> ]
```

Self-Attention

"Attention" is essentially a weighted correlation between inputs, of context size N and embedding (hidden) dimension d:

$$X_{N\times d} = [x_1, x_2, \dots, x_N], \text{ where } x_i \in \mathbb{R}^d$$

• Query (Q), Key (K) and Values (V) are linear transformations of X using weight matrices of dimensions $d \times d_k$ (typically $d_k = d$ for single-head attention).

$$Q_{N \times d_k} = XW_Q, \qquad K_{N \times d_k} = XW_K, \qquad V_{N \times d_k} = XW_V$$

In the attention weight, each cell represents how much one token attends to another. It's normalized so the weights sum to 1 across row.

$$Attention_{N \times N} = softmax \left(\frac{QK^T}{\sqrt{d_R}} \right)$$

The output is the attention weights multiplied by the values (weighted inputs):

$$Y_{N \times d_k} = \operatorname{softmax} \left(\frac{QK^T}{\sqrt{d_k}} \right) V = \operatorname{softmax} \left(\frac{X(W_Q W_K^T) X^T}{\sqrt{d_k}} \right) XW_V$$

Softmax

- The softmax (multinomial logistic) function is a multivariate version of the logistic function.
- It converts a vector z of K real numbers into a probability distribution on K categories.

$$\sigma_i(z) = \frac{e^{z_i}}{\sum_{i=1}^K e^{z_i}},$$
 where $i = 1, 2, \dots, K$

Note that it sums to one.

$$\sum_{i=1}^K \sigma_i =$$

In neural networks, softmax function is often used as the last activation function to convert network output into predicated output classes.

Simple Self-Attention (without Weights)

This replicates the codes in Section 3.3.

```
inputs <- torch::torch_tensor(rbind(c(0.43, 0.15, 0.89), # Your
                                  c(0.55, 0.87, 0.66), # journey
                                  c(0.57, 0.85, 0.64), # starts
                                  c(0.22, 0.58, 0.33), # with
                                  c(0.77, 0.25, 0.10), # one
                                  c(0.05, 0.80, 0.55))) # step - X is a 6 x 3 matrix
attn.score <- torch::torch mm(inputs.inputs$transpose(1, 2)) # A = X ** X ^T
## normalize by row (2nd dimension)
(attn.wghts <- torch::nn softmax(2)(attn.score))
                                                               #W = softmax(A), each row sums to 1
## torch tensor
## 0.2098 0.2006 0.1981 0.1242 0.1220 0.1452
## 0.1385 0.2379 0.2333 0.1240 0.1082 0.1581
## 0.1390 0.2369 0.2326 0.1242 0.1108 0.1565
## 0.1435 0.2074 0.2046 0.1462 0.1263 0.1720
## 0.1526 0.1958 0.1975 0.1367 0.1879 0.1295
## 0.1385 0.2184 0.2128 0.1420 0.0988 0.1896
## [ CPUFloatType{6.6} ]
(outputs <- torch::torch mm(attn.wghts. inputs)) # W %*% X
## torch tensor
## 0.4421 0.5931 0.5790
   0.4419 0.6515 0.5683
## 0.4431 0.6496 0.5671
## 0.4304 0.6208 0.5510
## 0.4671 0.5910 0.5266
## 0.4177 0.6503 0.5645
## [ CPUFloatType{6.3} ]
```

Self-Attention with Weights

This replicates the code in Section 3.4. The nn_linear function returns a *function* with optimized initial weights and performs the linear transformation (multiplication) efficiently.

```
d.in <- d.out <- ncol(inputs)
W.guery <- torch::nn linear(d.in, d.out, bias = FALSE)</pre>
W.key <- torch::nn linear(d.in. d.out. bias = FALSE)
W.value <- torch::nn linear(d.in, d.out, bias = FALSE)
self.attn.score <- torch::torch mm(W.query(inputs), W.key(inputs)$transpose(1, 2)) # As = X W Q %*% (X W K)^T
(self.attn.wghts <- torch::nn softmax(2)(self.attn.score / sgrt(d.out))) # Aw = softmax(A)
## torch tensor
## 0.1637 0.1623 0.1620 0.1737 0.1626 0.1757
## 0.1594 0.1630 0.1629 0.1743 0.1651 0.1752
## 0.1595 0.1633 0.1632 0.1739 0.1656 0.1745
## 0.1624 0.1644 0.1643 0.1715 0.1651 0.1724
## 0.1625 0.1686 0.1690 0.1646 0.1742 0.1611
## 0.1615 0.1619 0.1615 0.1757 0.1607 0.1789
## [ CPUFloatType{6.6} ][ grad fn = <SoftmaxBackwardo> ]
(outputs <- torch::torch mm(self.attn.wghts, W.value(inputs))) # Aw *** W V X
## torch tensor
## -0.3640 0.0766 0.2043
## -0.3647 0.0764 0.2053
## -0.3648 0.0761 0.2054
## -0.3648 0.0751 0.2052
## -0.3657 0.0689 0.2069
## -0.3643 0.0786 0.2044
## [ CPUFloatType{6,3} ][ grad fn = <MmBackwardo> ]
```

Causal Attention

 Casual (masked) attention restricts a model to only consider correlation with previous tokens, by multiplying the attention weights with a mask and re-normalize.

```
mask <- torch::torch_tril(torch::torch_ones(nrow(inputs), nrow(inputs)))</pre>
```

Alternatively, apply the mask first on the attention scores and re-normalize (-Inf is treated as o by softmax function).

```
(casual.attn.score <- self.attn.score$masked_fil(mask < 0.5, -Inf))
## torch_tensor
## -0.1961
             -inf
                     -inf
                             -inf
                                     -inf
                                             -inf
## -0.2356 -0.1969
                     -inf
                             -inf
                                     -inf
                                             -inf
                                             -inf
## -0.2259 -0.1853 -0.1868
                             -inf
                                     -inf
## -0.1446 -0.1233 -0.1248 -0.0503
                                     -inf
                                             -inf
## 0.0126 0.0762 0.0804 0.0341 0.1322
                                             -inf
## -0.2466 -0.2426 -0.2469 -0.1008 -0.2552 -0.0694
## [ CPUFloatType{6.6} ][ grad fn = <MaskedFillBackwardo> ]
(casual.attn.wghts <- torch::nn softmax(2)(casual.attn.score / sgrt(d.out)))</pre>
## torch_tensor
    1.0000 0.0000
                   0.0000
                           0.0000
                                   0.0000
                                           0.0000
   0.4944 0.5056 0.0000
                           0.0000 0.0000
                                           0.0000
   0.3282 0.3360 0.3357
                           0.0000 0.0000
                                           0.0000
   0.2451 0.2481 0.2479
                           0.2588 0.0000
                                           0.0000
   0.1938 0.2010 0.2015
                           0.1962 0.2076 0.0000
   0.1615 0.1619 0.1615 0.1757 0.1607 0.1789
    CPUFloatType{6.6} ][ grad fn = <SoftmaxBackwardo> ]
```

Dropout

- Dropout in deep learning is a technique where randomly selected hidden layer units are ignored during training, effectively "dropping" them out.
- This helps prevent overfitting by ensuring that a model does not become overly reliant on any specific set of hidden layer units.
- Dropout is only used during training and is disabled afterward.

```
dropout <- torch::nn dropout(p = 0.5) # drop 50%
dropout(casual.attn.wghts)
## torch_tensor
   2 0000 0 0000 0 0000
                          0.0000
                                  െ രെരെ
   0.9888 1.0112 0.0000
                          0.0000 0.0000
   0.6565 0.0000 0.6715
                          0.0000
                                  0.0000
                                         0.0000
   0.0000 0.0000 0.4959
                          0.0000
                                  0.0000
                                         0.0000
   0.0000 0.4020 0.0000
                          0.3923
                                  0.0000
                                         0.0000
   0.3230 0.0000 0.3220 0.0000 0.3214 0.3577
    CPUFloatType{6.6} ][ grad fn = <MulBackwardo> ]
```

■ Note that the remaining weights are scaled by a factor of 1/p.

Multi-Head Attention

- The term "multi-head" refers to dividing the attention mechanism into multiple "heads," each with its own set of (smaller) weight matrices (parameters), operating independently.
 - Computational efficiency: smaller weight matrices to deal with and can be optimized in batches
 - Specialization: each head is allowed to "learn" a different aspect of the sequence syntax, semantics, coreference, etc.
- In practical terms, this involves creating multiple instances of the self-attention mechanism (input of length d, output of length d_k , where commonly $d_k = d/H$) each with its own weight matrices (Q, K, V), and then combining the outputs into a vector of length $H \times d_k$.
- Or more efficiently, we create the query, key and value tensors of dimension $d \times (H \cdot d_k)$ through one matrix multiplication and then split each of them into a tensor of $H \times d \times d_k$ for use in subsequent calculations.

Example of Tensor Multiplication

Here is an example of tensor multiplication which is carried out on the last two dimensions. Multiplying a tensor of size $1 \times 2 \times 4 \times 2$ with another of size $1 \times 2 \times 2 \times 4$ yields a tensor of size $1 \times 2 \times 2 \times 2$.

```
aa
## torch tensor
## (1.1...) =
## 0.2745 0.8993 0.0772 0.4066
  0.2775 0.9268 0.1479 0.4545
## (1.2...) =
## 0.6584 0.0390 0.3565 0.2318
## 0.8573 0.7388 0.5331 0.9737
## [ CPUFloatType{1,2,2,4} ]
torch::torch_matmul(aa, aa$transpose(3, 4))
## torch_tensor
## (1.1...) =
  1.0554 1.1059
   1.1059 1.1644
## (1.2...) =
    0.6158 1.0090
   1.0090 2.5131
## [ CPUFloatType{1.2.2.2} ]
```

Layer Normalization

- Layer normalization is used to improve numerical stability by adjusting the activations (outputs) of a neural network layer to have a mean of o and a variance of 1.
- In GPT-2 and modern transformer architectures, layer normalization is typically applied before and after the multi-head attention module, and before the final output layer.

```
batch.ex <- torch::torch_randn(2, 5)
                                                                          # tensor of 2 x 5
layer.ex <- torch::nn_sequential(torch::nn_linear(5, 6), torch::nn_relu()) # one linear layer (5 x 6) + RELU activation function
(out.ex <- laver.ex(batch.ex))
                                                                          # 2 x 6 tensor
## torch tensor
## 0.1025 0.0176 1.6049 0.0503 0.0000 0.8693
## 0.0000 0.0073 1.0538 0.0000 0.0000 0.1109
## [ CPUFloatType{2.6} ][ grad fn = <ReluBackwardo> ]
(out.ex - out.ex$mean(-1, # operating on the last dimension
                     keepdim = TRUE # output a 2 x 1 tensor. instead of dropping to a vector of length 2
                     ))/(sqrt(out.ex$var(-1, keepdim = TRUE.
                                         unbiased = FALSE # use n as denominator instead of unbiased n-1, this was used in GPT-2
                                           # and is TensorFlow's default behavior. doesn't matter much when n is large
                                         ) + 1e-5)) # a small bias to avoid division by zero
## torch tensor
## -0.5613 -0.7022 1.9317 -0.6479 -0.7314 0.7111
## -0.5060 -0.4872 2.2240 -0.5060 -0.5060 -0.2188
## [ CPUFloatType{2.6} ][ grad fn = <DivBackwardo> ]
```

Activation Functions

- Activation functions introduce non-linearity into a neural network and are used after each layer. A different activation function (e.g., softmax) may be used after the final output layer.
- Common activation functions:

RELU (Rectified Linear Unit)
$$f(x) = \max(0, x)$$

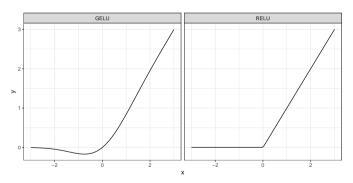
GELU (Gaussian error linear unit) $f(x) = x\Phi(x)$

$$\approx \frac{1}{2}x \left(1 + \tanh\left[\sqrt{\frac{2}{\pi}(x + 0.0044715 \cdot x^3)}\right]\right)$$

- $\Phi()$ is the CDF of normal distribution, and an approximation is used in the original GPT-2.
- The activation function is always applied element-wise to the input tensor. Note each element is often referred as a "neuron".

GELU vs RELU

■ GELU is similar to RELU but is a smooth function.



Shortcut Connection

- Shortcut connection adds the input of a layer (function) to its output (i.e., y = f(x) + x), which effectively adds 1 to the gradient (derivative) of the layer.
- It is designed to overcome the vanishing gradient problem, where the gradients become progressively smaller as they propagate backward through the layers, making it difficult to effectively train earlier layers.

Transform Block

- A feed forward layer is a small network that can be easily stacked to form more layers, because its input and output have the same dimension while at the same time allows more complexity in between.
 - 1. a linear layer that expands the embedding dimensions
 - 2. a nonlinear GELU activation function
 - 3. another linear layer that contracts the embedding dimensions to match the original input.
- A transformer block combines multi-head attention, layer normalization, dropout, feed forward layers, and GELU activations. It is the basic building block of a LLM and is repeated 12 times in the 124-million-parameter GPT-2.

Implementing a Transformer Block

```
x <- input
shortcut <- x
x <- layer_norm(x)</pre>
x <- multi head attention(x)</pre>
x <- dropout(x)
x <- x + shortcut
shortcut <- x
x <- laver norm(x)
x <- nn_sequential(nn_linear(d.in = emb.d, d.out = 4 * emb.d),
                    GELU().
                    nn linear(d.in = 4 * emb.d. d.out = emb.d))(x) # feed forward
x <- dropout(x)
output <- x + shortcut
layer_norm <- torch::nn_module(</pre>
    classname = "layer norm".
    initialize = function(emb.dim) {
        self$eps <- 1e-5
        self$scale <- torch::nn parameter(torch::torch ones(emb.dim))</pre>
        self$shift <- torch::nn parameter(torch::torch zeros(emb.dim))</pre>
    forward = function(x) {
        mean <- x$mean(dim = -1, keepdim = TRUE)
        var <- x$var(dim = -1. keepdim = TRUE, unbiased = FALSE)</pre>
        norm.x <- (x - mean) / torch::torch sqrt(var + self$eps)
        self$scale * norm.x + self$shift
ln <- laver norm(5)</pre>
```