```
In [ ]: #!pip install tensorflow==2.0.0
In [ ]: import tensorflow as tf
    print(tf.__version__)
2.3.0
```

Part 1: TensorFlow basics

Tensors

This is a constant tensor:

```
In []: x = tf.constant([[5, 2], [1, 3]])
        print(x)
       tf.Tensor(
       [[5 2]
        [1 \ 3], shape=(2, 2), dtype=int32)
        You can get its value as a Numpy array by calling .numpy():
In [ ]: x.numpy()
Out[]: array([[5, 2],
                [1, 3]], dtype=int32)
        Much like a Numpy array, it features the attributes dtype and shape:
In [ ]: print('dtype:', x.dtype)
        print('shape:', x.shape)
       dtype: <dtype: 'int32'>
       shape: (2, 2)
        A common way to create constant tensors is via tf.ones and tf.zeros (just
        like np.ones and np.zeros):
In [ ]: print(tf.ones(shape=(2, 1)))
        print(tf.zeros(shape=(2, 1)))
       tf.Tensor(
       [[1.]]
        [1.]], shape=(2, 1), dtype=float32)
       tf.Tensor(
       [[0]]
        [0.]], shape=(2, 1), dtype=float32)
```

Random constant tensors

This is all pretty normal:

Variables

Variables are special tensors used to store mutable state (like the weights of a neural network). You create a Variable using some initial value.

```
In [ ]: initial value = tf.random.normal(shape=(2, 2))
        a = tf.Variable(initial value)
        print(a)
       <tf.Variable 'Variable:0' shape=(2, 2) dtype=float32, numpy=
       array([[ 0.36464038, 1.8475662 ],
              [-0.5905565 , -0.16027513]], dtype=float32)>
        You update the value of a Variable by using the methods .assign(value), or
        .assign add(increment) or .assign sub(decrement):
In [ ]: new value = tf.random.normal(shape=(2, 2))
        a.assign(new value)
        for i in range(2):
          for j in range(2):
            assert a[i, j] == new value[i, j]
In []: added value = tf.random.normal(shape=(2, 2))
        a.assign add(added value)
        for i in range(2):
          for j in range(2):
            assert a[i, j] == new value[i, j] + added value[i, j]
```

Doing math in TensorFlow

You can use TensorFlow exactly like you would use Numpy. The main difference is that your TensorFlow code can run on GPU and TPU.

```
In []: a = tf.random.normal(shape=(2, 2))
b = tf.random.normal(shape=(2, 2))

c = a + b
d = tf.square(c)
e = tf.exp(d)
```

Computing gradients with GradientTape

Oh, and there's another big difference with Numpy: you can automatically retrieve the gradient of any differentiable expression.

Just open a GradientTape, start "watching" a tensor via tape.watch(), and compose a differentiable expression using this tensor as input:

```
In []: a = tf.random.normal(shape=(2, 2))
b = tf.random.normal(shape=(2, 2))

with tf.GradientTape() as tape:
    tape.watch(a) # Start recording the history of operations applied to `a`
    c = tf.sqrt(tf.square(a) + tf.square(b)) # Do some math using `a`
    # What's the gradient of `c` with respect to `a`?
    dc_da = tape.gradient(c, a)
    print(dc_da)

tf.Tensor(
[[ 0.80471563 -0.98697984]
    [-0.7596037    0.15738489]], shape=(2, 2), dtype=float32)
```

By default, variables are watched automatically, so you don't need to manually watch them:

```
In []: a = tf.Variable(a)

with tf.GradientTape() as tape:
    c = tf.sqrt(tf.square(a) + tf.square(b))
    dc_da = tape.gradient(c, a)
    print(dc_da)

tf.Tensor(
  [[ 0.80471563 -0.98697984]
  [-0.7596037   0.15738489]], shape=(2, 2), dtype=float32)
```

Note that you can compute higher-order derivatives by nesting tapes:

```
In []: with tf.GradientTape() as outer_tape:
    with tf.GradientTape() as tape:
        c = tf.sqrt(tf.square(a) + tf.square(b))
        dc_da = tape.gradient(c, a)
        d2c_da2 = outer_tape.gradient(dc_da, a)
        print(d2c_da2)
```

```
tf.Tensor(
[[0.176198    0.02943563]
  [0.3462551   1.219587 ]], shape=(2, 2), dtype=float32)
```

An end-to-end example: linear regression

So far you've learned that TensorFlow is a Numpy-like library that is GPU or TPU accelerated, with automatic differentiation. Time for an end-to-end example: let's implement a linear regression, the FizzBuzz of Machine Learning.

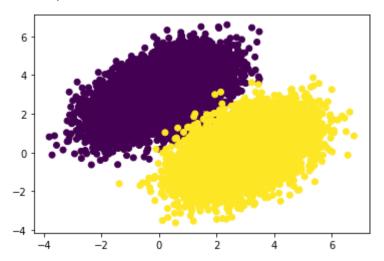
For the sake of demonstration, we won't use any of the higher-level Keras components like Layer or MeanSquaredError. Just basic ops.

```
In [ ]: input dim = 2
        output dim = 1
        learning rate = 0.01
        # This is our weight matrix
        w = tf.Variable(tf.random.uniform(shape=(input dim, output dim)))
        # This is our bias vector
        b = tf.Variable(tf.zeros(shape=(output dim,)))
        def compute predictions(features):
          return tf.matmul(features, w) + b
        def compute loss(labels, predictions):
          return tf.reduce mean(tf.square(labels - predictions))
        def train on batch(x, y):
          with tf.GradientTape() as tape:
            predictions = compute predictions(x)
            loss = compute loss(y, predictions)
            dloss dw, dloss db = tape.gradient(loss, [w, b])
          w.assign sub(learning rate * dloss dw)
          b.assign sub(learning rate * dloss db)
          return loss
```

Let's generate some artificial data to demonstrate our model:

```
plt.scatter(features[:, 0], features[:, 1], c=labels[:, 0])
```

Out[]: <matplotlib.collections.PathCollection at 0x7f29d00f5908>

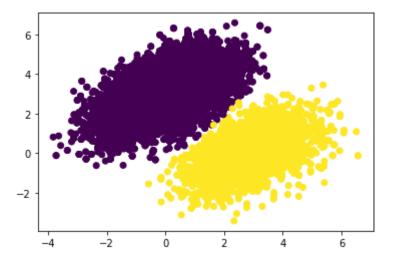


Now let's train our linear regression by iterating over batch-by-batch over the data and repeatedly calling train on batch:

```
In [ ]: # Shuffle the data.
        random.Random(1337).shuffle(features)
        random.Random(1337).shuffle(labels)
        # Create a tf.data.Dataset object for easy batched iteration
        dataset = tf.data.Dataset.from tensor slices((features, labels))
        dataset = dataset.shuffle(buffer size=1024).batch(256)
        for epoch in range(10):
          for step, (x, y) in enumerate(dataset):
            loss = train on batch(x, y)
          print('Epoch %d: last batch loss = %.4f' % (epoch, float(loss)))
       Epoch 0: last batch loss = 0.1011
       Epoch 1: last batch loss = 0.0772
       Epoch 2: last batch loss = 0.0601
       Epoch 3: last batch loss = 0.0427
       Epoch 4: last batch loss = 0.0352
       Epoch 5: last batch loss = 0.0388
       Epoch 6: last batch loss = 0.0369
       Epoch 7: last batch loss = 0.0300
       Epoch 8: last batch loss = 0.0234
       Epoch 9: last batch loss = 0.0266
        Here's how our model performs:
```

plt.scatter(features[:, 0], features[:, 1], c=predictions[:, 0] > 0.5)

In []: predictions = compute predictions(features)



Making it fast with tf.function

But how fast is our current code running?

```
In []: import time

t0 = time.time()
for epoch in range(20):
    for step, (x, y) in enumerate(dataset):
        loss = train_on_batch(x, y)
    t_end = time.time() - t0
    print('Time per epoch: %.3f s' % (t_end / 20,))
```

Time per epoch: 0.149 s

Let's compile the training function into a static graph. Literally all we need to do is add the tf.function decorator on it:

```
In []: @tf.function
    def train_on_batch(x, y):
        with tf.GradientTape() as tape:
            predictions = compute_predictions(x)
            loss = compute_loss(y, predictions)
            dloss_dw, dloss_db = tape.gradient(loss, [w, b])
        w.assign_sub(learning_rate * dloss_dw)
        b.assign_sub(learning_rate * dloss_db)
        return loss
```

Let's try this again:

```
In []: t0 = time.time()
for epoch in range(20):
    for step, (x, y) in enumerate(dataset):
        loss = train_on_batch(x, y)
    t_end = time.time() - t0
    print('Time per epoch: %.3f s' % (t_end / 20,))
```

Time per epoch: 0.084 s

40% reduction, neat. In this case we used a trivially simple model; in general the bigger the model the greater the speedup you can get by leveraging static graphs.

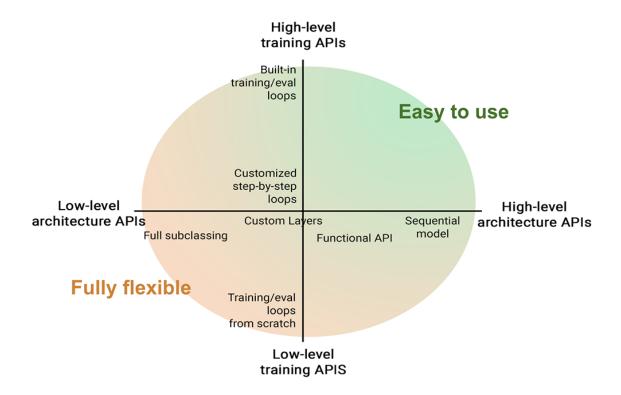
Remember: eager execution is great for debugging and printing results line-byline, but when it's time to scale, static graphs are a researcher's best friends.

Part 2: The Keras API

Keras is a Python API for deep learning. It has something for everyone:

- If you're an engineer, Keras provides you with reusable blocks such as layers, metrics, training loops, to support common use cases. It provides a high-level user experience that's accessible and productive.
- If you're a researcher, you may prefer not to use these built-in blocks such as
 layers and training loops, and instead create your own. Of course, Keras
 allows you to do this. In this case, Keras provides you with templates for the
 blocks you write, it provides you with structure, with an API standard for
 things like Layers and Metrics. This structure makes your code easy to share
 with others and easy to integrate in production workflows.
- The same is true for library developers: TensorFlow is a large ecosystem. It
 has many different libraries. In order for different libraries to be able to talk
 to each other and share components, they need to follow an API standard.
 That's what Keras provides.

Crucially, Keras brings high-level UX and low-level flexibility together fluently: you no longer have on one hand, a high-level API that's easy to use but inflexible, and on the other hand a low-level API that's flexible but only approachable by experts. Instead, you have a spectrum of workflows, from the very high-level to the very low-level. Workflows that are all compatible because they're built on top of the same concepts and objects.



The base Layer class

The first class you need to know is Layer . Pretty much everything in Keras derives from it.

A Layer encapsulates a state (weights) and some computation (defined in the call method).

```
In [ ]: from tensorflow.keras.layers import Layer
        class Linear(Layer):
          """v = w.x + b"""
          def init (self, units=32, input dim=32):
              super(Linear, self).__init__()
              w init = tf.random normal initializer()
              self.w = tf.Variable(
                  initial value=w init(shape=(input dim, units), dtype='float32'),
                  trainable=True)
              b init = tf.zeros initializer()
              self.b = tf.Variable(
                  initial value=b init(shape=(units,), dtype='float32'),
                  trainable=True)
          def call(self, inputs):
              return tf.matmul(inputs, self.w) + self.b
        # Instantiate our layer.
        linear layer = Linear(4, 2)
```

A layer instance works like a function. Let's call it on some data:

w init = tf.random normal initializer()

us from having to specify input dim in the constructor:

```
In [ ]: y = linear_layer(tf.ones((2, 2)))
assert y.shape == (2, 4)
```

The Layer class takes care of tracking the weights assigned to it as attributes:

```
In [ ]: # Weights are automatically tracked under the `weights` property.
assert linear_layer.weights == [linear_layer.w, linear_layer.b]
```

Note that's also a shortcut method for creating weights: <code>add_weight</code> . Instead of doing

```
self.w = tf.Variable(initial_value=w_init(shape=shape,
dtype='float32'))
You would typically do:
self.w = self.add weight(shape=shape, initializer='random normal')
```

It's good practice to create weights in a separate build method, called lazily with the shape of the first inputs seen by your layer. Here, this pattern prevents

```
In []: class Linear(Layer):
    """y = w.x + b"""

    def __init__(self, units=32):
        super(Linear, self).__init__()

Loading [MathJax]/jax/output/CommonHTML/fonts/TeX/fontdata.js
```

Trainable and non-trainable weights

Weights created by layers can be either trainable or non-trainable. They're exposed in trainable_weights and non_trainable_weights. Here's a layer with a non-trainable weight:

```
In [ ]: class ComputeSum(Layer):
          """Returns the sum of the inputs."""
          def init (self, input dim):
              super(ComputeSum, self).__init__()
              # Create a non-trainable weight.
              self.total = tf.Variable(initial value=tf.zeros((input dim,)),
                                       trainable=False)
          def call(self, inputs):
              self.total.assign add(tf.reduce sum(inputs, axis=0))
              return self.total
        my sum = ComputeSum(2)
        x = tf.ones((2, 2))
        y = my sum(x)
        print(y.numpy()) # [2. 2.]
        y = my sum(x)
        print(y.numpy()) # [4. 4.]
        assert my sum.weights == [my sum.total]
        assert my sum.non trainable weights == [my sum.total]
        assert my sum.trainable weights == []
```

Recursively composing layers

Layers can be recursively nested to create bigger computation blocks. Each layer will track the weights of its sublayers (both trainable and non-trainable.

```
In [ ]: # Let's reuse the Linear class
        # with a `build` method that we defined above.
        class MLP(Layer):
            """Simple stack of Linear layers."""
            def init (self):
                super(MLP, self).__init__()
                self.linear 1 = Linear(32)
                self.linear 2 = Linear(32)
                self.linear_3 = Linear(10)
            def call(self, inputs):
                x = self.linear 1(inputs)
                x = tf.nn.relu(x)
                x = self.linear 2(x)
                x = tf.nn.relu(x)
                return self.linear 3(x)
        mlp = MLP()
        # The first call to the `mlp` object will create the weights.
        y = mlp(tf.ones(shape=(3, 64)))
        # Weights are recursively tracked.
        assert len(mlp.weights) == 6
```

Built-in layers

Keras provides you with a wide range of built-in layers, so that you don't have to implement your own layers all the time.

- Convolution layers
- Transposed convolutions
- Separateable convolutions
- Average and max pooling
- Global average and max pooling
- LSTM, GRU (with built-in cuDNN acceleration)
- BatchNormalization
- Dropout
- Attention

- ConvLSTM2D
- etc.

Keras follows the principles of exposing good default configurations, so that layers will work fine out of the box for most use cases if you leave keyword arguments to their default value. For instance, the LSTM layer uses an orthogonal recurrent matrix initializer by default, and initializes the forget gate bias to one by default.

The training argument in call

Some layers, in particular the BatchNormalization layer and the Dropout layer, have different behaviors during training and inference. For such layers, it is standard practice to expose a training (boolean) argument in the call method.

By exposing this argument in call, you enable the built-in training and evaluation loops (e.g. fit) to correctly use the layer in training and inference.

```
In [ ]: class Dropout(Layer):
          def init (self, rate):
            super(Dropout, self).__init__()
            self.rate = rate
          def call(self, inputs, training=None):
            if training:
              return tf.nn.dropout(inputs, rate=self.rate)
            return inputs
        class MLPWithDropout(Layer):
          def init (self):
              super(MLPWithDropout, self). init ()
              self.linear 1 = Linear(32)
              self.dropout = Dropout(0.5)
              self.linear 3 = Linear(10)
          def call(self, inputs, training=None):
              x = self.linear 1(inputs)
              x = tf.nn.relu(x)
              x = self.dropout(x, training=training)
              return self.linear 3(x)
        mlp = MLPWithDropout()
        y train = mlp(tf.ones((2, 2)), training=True)
        y test = mlp(tf.ones((2, 2)), training=False)
```

To build deep learning models, you don't have to use object-oriented programming all the time. Layers can also be composed functionally, like this (we call it the "Functional API"):

```
In [ ]: # We use an `Input` object to describe the shape and dtype of the inputs.
        # This is the deep learning equivalent of *declaring a type*.
        # The shape argument is per-sample; it does not include the batch size.
        # The functional API focused on defining per-sample transformations.
        # The model we create will automatically batch the per-sample transformation
        # so that it can be called on batches of data.
        inputs = tf.keras.Input(shape=(16,))
        # We call layers on these "type" objects
        # and they return updated types (new shapes/dtypes).
        x = Linear(32)(inputs) # We are reusing the Linear layer we defined earlier.
        x = Dropout(0.5)(x) # We are reusing the Dropout layer we defined earlier.
        outputs = Linear(10)(x)
        # A functional `Model` can be defined by specifying inputs and outputs.
        # A model is itself a layer like any other.
        model = tf.keras.Model(inputs, outputs)
        # A functional model already has weights, before being called on any data.
        # That's because we defined its input shape in advance (in `Input`).
        assert len(model.weights) == 4
        # Let's call our model on some data.
        y = model(tf.ones((2, 16)))
        assert y.shape == (2, 10)
```

The Functional API tends to be more concise than subclassing, and provides a few other advantages (generally the same advantages that functional, typed languages provide over untyped OO development). However, it can only be used to define DAGs of layers -- recursive networks should be defined as Layer subclasses instead.

Key differences between models defined via subclassing and Functional models are explained in this blog post.

Learn more about the Functional API here.

In your research workflows, you may often find yourself mix-and-matching OO models and Functional models.

For models that are simple stacks of layers with a single input and a single output, you can also use the Sequential class which turns a list of layers into a Model:

```
model = Sequential([Linear(32), Dropout(0.5), Linear(10)])

y = model(tf.ones((2, 16)))
assert y.shape == (2, 10)
```

Loss classes

Keras features a wide range of built-in loss classes, like BinaryCrossentropy, CategoricalCrossentropy, KLDivergence, etc. They work like this:

```
In [ ]: bce = tf.keras.losses.BinaryCrossentropy()
    y_true = [0., 0., 1., 1.] # Targets
    y_pred = [1., 1., 1., 0.] # Predictions
    loss = bce(y_true, y_pred)
    print('Loss:', loss.numpy())
```

Loss: 11.522857

Note that loss classes are stateless: the output of __call__ is only a function of the input.

Metric classes

Keras also features a wide range of built-in metric classes, such as BinaryAccuracy, AUC, FalsePositives, etc.

Unlike losses, metrics are stateful. You update their state using the update_state method, and you query the scalar metric result using result:

```
In []: m = tf.keras.metrics.AUC()
    m.update_state([0, 1, 1, 1], [0, 1, 0, 0])
    print('Intermediate result: ', m.result().numpy())

m.update_state([1, 1, 1, 1], [0, 1, 1, 0])
    print('Final result: ', m.result().numpy())
```

Intermediate result: 0.6666667 Final result: 0.71428573

The internal state can be cleared with metric.reset_states.

You can easily roll out your own metrics by subclassing the Metric class:

- Create the state variables in init
- Update the variables given y true and y pred in update state
- Return the metric result in result
- Clear the state in reset states

Here's a quick implementation of a BinaryTruePositives metric as a demonstration:

```
In [ ]: class BinaryTruePositives(tf.keras.metrics.Metric):
          def init (self, name='binary true positives', **kwargs):
            super(BinaryTruePositives, self). init (name=name, **kwargs)
            self.true positives = self.add weight(name='tp', initializer='zeros')
          def update_state(self, y_true, y_pred, sample_weight=None):
            y true = tf.cast(y true, tf.bool)
            y pred = tf.cast(y pred, tf.bool)
            values = tf.logical and(tf.equal(y true, True), tf.equal(y pred, True))
            values = tf.cast(values, self.dtype)
            if sample weight is not None:
              sample weight = tf.cast(sample weight, self.dtype)
              sample weight = tf.broadcast weights(sample weight, values)
              values = tf.multiply(values, sample weight)
            self.true positives.assign add(tf.reduce sum(values))
          def result(self):
            return self.true positives
          def reset states(self):
            self.true positive.assign(0)
```

Optimizer classes & a quick end-to-end training loop

You don't normally have to define by hand how to update your variables during gradient descent, like we did in our initial linear regression example. You would usually use one of the built-in Keras optimizer, like SGD, RMSprop, or Adam.

Here's a simple MNSIT example that brings together loss classes, metric classes, and optimizers.

```
# Instantiate a logistic loss function that expects integer targets.
loss = tf.keras.losses.SparseCategoricalCrossentropy(from logits=True)
# Instantiate an accuracy metric.
accuracy = tf.keras.metrics.SparseCategoricalAccuracy()
# Instantiate an optimizer.
optimizer = tf.keras.optimizers.Adam()
# Iterate over the batches of the dataset.
for step, (x, y) in enumerate(dataset):
 # Open a GradientTape.
 with tf.GradientTape() as tape:
    # Forward pass.
   logits = model(x)
    # Loss value for this batch.
   loss value = loss(y, logits)
 # Get gradients of weights wrt the loss.
 gradients = tape.gradient(loss value, model.trainable weights)
 # Update the weights of our linear layer.
 optimizer.apply gradients(zip(gradients, model.trainable weights))
 # Update the running accuracy.
 accuracy.update state(y, logits)
 # Logging.
 if step % 100 == 0:
   print('Step:', step)
    print('Loss from last step:', float(loss_value))
    print('Total running accuracy so far:', float(accuracy.result()))
```

```
Downloading data from https://storage.googleapis.com/tensorflow/tf-keras-dat
asets/mnist.npz
Step: 0
Loss from last step: 2.358793258666992
Total running accuracy so far: 0.09375
Step: 100
Loss from last step: 0.21707287430763245
Total running accuracy so far: 0.8310643434524536
Step: 200
Loss from last step: 0.2818300127983093
Total running accuracy so far: 0.8765547275543213
Step: 300
Loss from last step: 0.23447920382022858
Total running accuracy so far: 0.8955564498901367
Step: 400
Loss from last step: 0.11367885768413544
Total running accuracy so far: 0.9080813527107239
Step: 500
Loss from last step: 0.11368697881698608
Total running accuracy so far: 0.9158245921134949
Step: 600
Loss from last step: 0.0994415432214737
Total running accuracy so far: 0.9220309853553772
Step: 700
Loss from last step: 0.047019436955451965
Total running accuracy so far: 0.9272022247314453
Step: 800
Loss from last step: 0.07821480929851532
Total running accuracy so far: 0.9307116270065308
Step: 900
Loss from last step: 0.09753896296024323
Total running accuracy so far: 0.9345518946647644
```

We can reuse our SparseCategoricalAccuracy metric instance to implement a testing loop:

```
In []: x_test = x_test[:].reshape(10000, 784).astype('float32') / 255
    test_dataset = tf.data.Dataset.from_tensor_slices((x_test, y_test))
    test_dataset = test_dataset.batch(128)

accuracy.reset_states() # This clears the internal state of the metric

for step, (x, y) in enumerate(test_dataset):
    logits = model(x)
    accuracy.update_state(y, logits)

print('Final test accuracy:', float(accuracy.result()))
```

Final test accuracy: 0.9607999920845032

The add_loss method

Sometimes you need to compute loss values on the fly during a foward pass (especially regularization losses). Keras allows you to compute loss values at any time, and to recursively keep track of them via the add loss method.

Here's an example of a layer that adds a sparsity regularization loss based on the L2 norm of the inputs:

```
In []:
    class ActivityRegularization(Layer):
        """Layer that creates an activity sparsity regularization loss."""

    def __init__(self, rate=le-2):
        super(ActivityRegularization, self).__init__()
        self.rate = rate

    def call(self, inputs):
        # We use `add_loss` to create a regularization loss
        # that depends on the inputs.
        self.add_loss(self.rate * tf.reduce_sum(tf.square(inputs)))
        return inputs
```

Loss values added via add_loss can be retrieved in the .losses list property of any Layer or Model:

```
In []: from tensorflow.keras import layers

class SparseMLP(Layer):
    """Stack of Linear layers with a sparsity regularization loss."""

def __init__(self, output_dim):
    super(SparseMLP, self).__init__()
    self.dense_1 = layers.Dense(32, activation=tf.nn.relu)
    self.regularization = ActivityRegularization(1e-2)
    self.dense_2 = layers.Dense(output_dim)

def call(self, inputs):
    x = self.dense_1(inputs)
    x = self.regularization(x)
    return self.dense_2(x)

mlp = SparseMLP(1)
    y = mlp(tf.ones((10, 10)))

print(mlp.losses) # List containing one float32 scalar
```

[<tf.Tensor: shape=(), dtype=float32, numpy=0.6885692>]

These losses are cleared by the top-level layer at the start of each forward pass - they don't accumulate. So layer.losses always contain only the losses created during the last forward pass. You would typically use these losses by summing them before computing your gradients when writing a training loop.

```
In [ ]: # Losses correspond to the *last* forward pass.
        mlp = SparseMLP(1)
        mlp(tf.ones((10, 10)))
        assert len(mlp.losses) == 1
        mlp(tf.ones((10, 10)))
        assert len(mlp.losses) == 1 # No accumulation.
        # Let's demonstrate how to use these losses in a training loop.
        # Prepare a dataset.
        (x train, y train), = tf.keras.datasets.mnist.load data()
        dataset = tf.data.Dataset.from tensor slices(
            (x train.reshape(60000, 784).astype('float32') / 255, y train))
        dataset = dataset.shuffle(buffer size=1024).batch(64)
        # A new MLP.
        mlp = SparseMLP(10)
        # Loss and optimizer.
        loss fn = tf.keras.losses.SparseCategoricalCrossentropy(from logits=True)
        optimizer = tf.keras.optimizers.SGD(learning rate=0.1)
        for step, (x, y) in enumerate(dataset):
          with tf.GradientTape() as tape:
            # Forward pass.
            logits = mlp(x)
            # External loss value for this batch.
            loss = loss fn(y, logits)
            # Add the losses created during the forward pass.
            loss += sum(mlp.losses)
            # Get gradients of weights wrt the loss.
            gradients = tape.gradient(loss, mlp.trainable weights)
          # Update the weights of our linear layer.
          optimizer.apply gradients(zip(gradients, mlp.trainable weights))
          # Logging.
          if step % 100 == 0:
            print(step, float(loss))
       0 4.304479598999023
       100 2.2919511795043945
       200 2.2856969833374023
       300 2.249835968017578
       400 2.154803514480591
       500 2.179860830307007
       600 2.0276057720184326
       700 2.064443349838257
       800 2.0333380699157715
       900 1.828589916229248
```

A detailed end-to-end example: a Variational AutoEncoder (VAE)

If you want to take a break from the basics and look at a slightly more advanced example, check out this Variational AutoEncoder implementation that demonstrates everything you've learned so far:

- Subclassing Layer
- Recursive layer composition
- Loss classes and metric classes
- add loss
- GradientTape

Using built-in training loops

It would be a bit silly if you had to write your own low-level training loops every time for simple use cases. Keras provides you with a built-in training loop on the Model class. If you want to use it, either subclass from Model or create a Functional or Sequential model.

To demonstrate it, let's reuse the MNIST setup from above:

```
In [ ]: # Prepare a dataset.
        (x_train, y_train), (x_test, y_test) = tf.keras.datasets.mnist.load_data()
        x train = x train.reshape(60000, 784).astype('float32') / 255
        dataset = tf.data.Dataset.from tensor slices((x train, y train))
        dataset = dataset.shuffle(buffer size=1024).batch(64)
        # Instantiate a simple classification model
        model = tf.keras.Sequential([
          layers.Dense(256, activation=tf.nn.relu),
          layers.Dense(256, activation=tf.nn.relu),
          layers.Dense(10)
        ])
        # Instantiate a logistic loss function that expects integer targets.
        loss = tf.keras.losses.SparseCategoricalCrossentropy(from logits=True)
        # Instantiate an accuracy metric.
        accuracy = tf.keras.metrics.SparseCategoricalAccuracy()
        # Instantiate an optimizer.
        optimizer = tf.keras.optimizers.Adam()
```

First, call compile to configure the optimizer, loss, and metrics to monitor.

Then we call fit on our model to pass it the data:

```
In [ ]: model.fit(dataset, epochs=3)
         Epoch 1/3
         se categorical accuracy: 0.9352
         Epoch 2/3
         se categorical accuracy: 0.9737
         Epoch 3/3
         se categorical accuracy: 0.9817
   Out[]: <tensorflow.python.keras.callbacks.History at 0x7f29713263c8>
          Done! Now let's test it:
   In []: x \text{ test} = x \text{ test}[:].reshape(10000, 784).astype('float32') / 255
          test dataset = tf.data.Dataset.from tensor slices((x test, y test))
          test dataset = test dataset.batch(128)
          loss, acc = model.evaluate(test dataset)
          print('loss:', loss, 'acc:', acc)
         79/79 [============= ] - Os 2ms/step - loss: 0.0848 - sparse
          categorical accuracy: 0.9747
         loss: 0.08479571342468262 acc: 0.9746999740600586
          Note that you can also monitor your loss and metrics on some validation data
          during fit.
          Also, you can call fit directly on Numpy arrays, so no need for the dataset
          conversion:
   In [ ]: (x train, y train), (x test, y test) = tf.keras.datasets.mnist.load data()
          x train = x train.reshape(60000, 784).astype('float32') / 255
          num val samples = 10000
          x val = x train[-num val samples:]
          y val = y train[-num val samples:]
          x train = x train[:-num val samples]
          y train = y train[:-num val samples]
          # Instantiate a simple classification model
          model = tf.keras.Sequential([
            layers.Dense(256, activation=tf.nn.relu),
            layers.Dense(256, activation=tf.nn.relu),
            layers.Dense(10)
          ])
          # Instantiate a logistic loss function that expects integer targets.
          loss = tf.keras.losses.SparseCategoricalCrossentropy(from logits=True)
Loading [MathJax]/jax/output/CommonHTML/fonts/TeX/fontdata.js | metric.
```

```
accuracy = tf.keras.metrics.SparseCategoricalAccuracy()
# Instantiate an optimizer.
optimizer = tf.keras.optimizers.Adam()
model.compile(optimizer=optimizer,
              loss=loss,
              metrics=[accuracy])
model.fit(x train, y train,
          validation data=(x val, y val),
          epochs=3,
          batch size=64)
```

```
Epoch 1/3
    se categorical accuracy: 0.9276 - val loss: 0.1197 - val sparse categorical
    accuracy: 0.9645
    Epoch 2/3
    se categorical accuracy: 0.9714 - val loss: 0.0982 - val sparse categorical
    accuracy: 0.9723
    Epoch 3/3
    se categorical accuracy: 0.9810 - val loss: 0.0774 - val sparse categorical
    accuracy: 0.9763
Out[]: <tensorflow.python.keras.callbacks.History at 0x7f2970435860>
```

Callbacks

One of the neat features of fit (besides built-in support for sample weighting and class weighting) is that you can easily customize what happens during training and evaluation by using callbacks.

A callback is an object that is called at different points during training (e.g. at the end of every batch or at the end of every epoch) and does stuff.

There's a bunch of built-in callback available, like ModelCheckpoint to save your models after each epoch during training, or EarlyStopping, which interrupts training when your validation metrics start stalling.

And you can easily write your own callbacks.

```
In [ ]: # Instantiate a simple classification model
        model = tf.keras.Sequential([
          layers.Dense(256, activation=tf.nn.relu),
          layers.Dense(256, activation=tf.nn.relu),
          layers.Dense(10)
        1)
        # Instantiate a logistic loss function that expects integer targets.
        loss = tf.keras.losses.SparseCategoricalCrossentropy(from logits=True)
```

```
# Instantiate an accuracy metric.
 accuracy = tf.keras.metrics.SparseCategoricalAccuracy()
 # Instantiate an optimizer.
 optimizer = tf.keras.optimizers.Adam()
model.compile(optimizer=optimizer,
           loss=loss,
           metrics=[accuracy])
 # Instantiate some callbacks
 callbacks = [tf.keras.callbacks.EarlyStopping(),
          tf.keras.callbacks.ModelCheckpoint(filepath='my model.keras',
                                      save best only=True)]
model.fit(x train, y train,
        validation data=(x val, y val),
        epochs=30,
        batch size=64,
        callbacks=callbacks)
Epoch 1/30
se categorical accuracy: 0.9287 - val loss: 0.1373 - val sparse categorical
accuracy: 0.9582
Epoch 2/30
782/782 [============] - 2s 2ms/step - loss: 0.0949 - spar
se categorical accuracy: 0.9706 - val loss: 0.0981 - val sparse categorical
accuracy: 0.9711
Epoch 3/30
se categorical accuracy: 0.9811 - val loss: 0.0944 - val sparse categorical
accuracy: 0.9717
Epoch 4/30
se categorical accuracy: 0.9851 - val loss: 0.0802 - val sparse categorical
accuracy: 0.9768
Epoch 5/30
se categorical accuracy: 0.9893 - val loss: 0.0800 - val sparse categorical
accuracy: 0.9780
```

Out[]: <tensorflow.python.keras.callbacks.History at 0x7f296e2fee10>

Parting words

I hope this guide has given you a good overview of what's possible with TensorFlow 2.0 and Keras!

Epoch 6/30

accuracy: 0.9755

Remember that TensorFlow and Keras don't represent a single workflow. It's a spectrum of workflows, each with its own trade-off between usability and flexibility. For instance, you've noticed that it's much easier to use fit than to write a custom training loop, but fit doesn't give you the same level of granular control for research use cases.

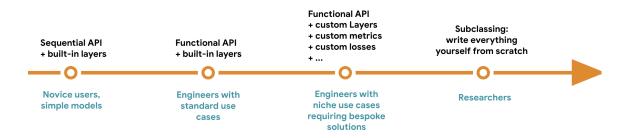
So use the right tool for the job!

A core principle of Keras is "progressive disclosure of complexity": it's easy to get started, and you can gradually dive into workflows where you write more and more logic from scratch, providing you with complete control.

This applies to both model definition, and model training.

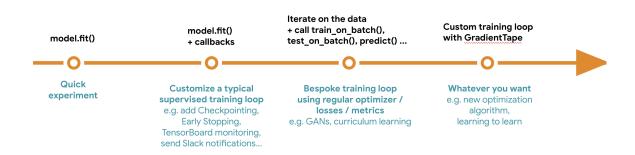
Model building: from simple to arbitrarily flexible

Progressive disclosure of complexity



Model training: from simple to arbitrarily flexible

Progressive disclosure of complexity



What to learn next

- Saving and serialization
- Distributed training on multiple GPUS
- Exporting models to TFLite for deployment on Android or embedded systems
- Exporting models to TensorFlow.js for deployment in the browser