ml-pipeline-refactoring

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0.1 Machine Learning Engineer @ InstaDeep - Technical test

What is this all about? This notebook provides a code snippet to implement a protein classifier. Specifically, for each protein, the task is to assign the corresponding Pfam family (i.e. protein family). You can find more information regarding the Pfam family here.

The implemented model is inspired by ProtCNN.

Where does the data come from? The PFAM dataset is available here.

What we are expecting? While this notebook provides a baseline model and some util functions to train the latter, its architecture is not suited for reproducibility, collaboration and portability. Therefore, your task is to refactor and augment the following basecode. In practice, we expect you to create a git repository and apply your suggestions, it should include but not limited:

- Build a local environment using docker
- Clean & refactor the code into python scripts (specifically, the user should be able to tweak training & model hyperparameters from the CLI)
- Add a command line to predict and/or evaluate a trained model
- Ensure code quality & consistency
- Document the repository
- Add tests
- ...

The above list is not exhaustive and we are more than open to other suggestions and inititatives.

- Bonus: Find a stronger model: e.g. optimize over hyperparameters, refine or change the architecture, etc.
- If you use GitHub/GitLab to host your git repository, please make it **private**. To share your work with us, please launch from your repository git bundle create <YOUR_NAME>.bundle --all and send the resulting <YOUR_NAME>.bundle file to us.
- The notebook has been tested with python==3.7.4 and the following dependencies:

```
matplotlib==3.4.1
numpy==1.18.5
pandas==1.2.3
pytorch-lightning==1.5.3
seaborn==0.11.1
```

```
tensorboard==2.2.2
torch==1.8.1
torchmetrics==0.6.0
```

- We use PyTorch Lightning for this pipeline's deep learning framework, but you are allowed to switch to any other ML/DL framework that you might prefer.
- Warning: The default number of GPUs is set to 0 but training without a GPU will be slow. In case you need more compute power than locally available on your computer, the following resources provide interesting amounts of computing power for free:
 - Google Colab: access to one GPU or one TPU, time limit of 12 hours (kernels are shut down after 12 hours)
 - Kaggle notebooks: access to one GPU (NVIDIA P100), time limit of 6 hours

```
[1]: import os

from collections import Counter

import pandas as pd
import numpy as np

import matplotlib.pyplot as plt
import seaborn as sns

import torch
import torch
import pytorch_lightning as pl
import torchmetrics
```

/Users/theomeborckinstadeep/Desktop/experiments/venv/lib/python3.7/site-packages/pandas/compat/__init__.py:97: UserWarning: Could not import the lzma module. Your installed Python is incomplete. Attempting to use lzma compression will result in a RuntimeError.

warnings.warn(msg)

```
[2]: torch.cuda.is_available()
```

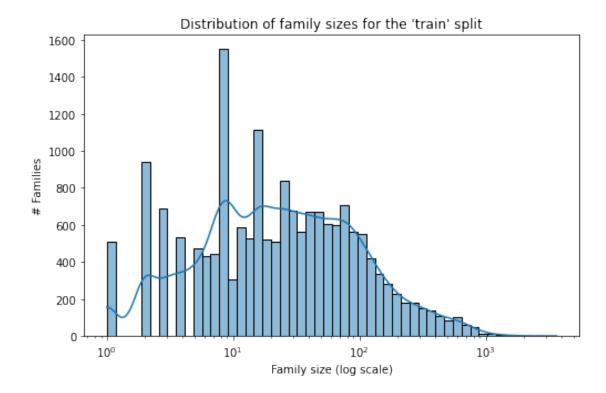
[2]: False

```
[3]: data_dir = './random_split'
```

0.1.1 Step 1: Define data loading and pre-processing

```
all_data = pd.concat(data)
         return all_data["sequence"], all_data["family_accession"]
[5]: train_data, train_targets = reader("train", data_dir)
[6]: train_data.head()
[6]: 0
          ISEAELEIMKVLWLKSPQTANEIIEELEDPMDWKPKTIRTLINRLV...
     1
          AGVPCSVKASEGYLFPLDRCFLFVTKPTLYIPYSEISSVVMSRTGG...
          IRHVLMNSPPGKLYDLVKDINILLGSSVSIQKILEEVLKDYNEKNY...
         MCIAIPGRIERIDYPIAIVDFKGLKKEVRIDLLENPQIGDYVLVHV...
          NIFHILWEDVDLEGVTFKPMGESISVQGDIHIFVLYEGEGENTPIR...
     Name: sequence, dtype: object
[7]: def build_labels(targets):
         unique_targets = targets.unique()
         fam2label = {target: i for i, target in enumerate(unique_targets, start=1)}
         fam2label['<unk>'] = 0
         print(f"There are {len(fam2label)} labels.")
         return fam2label
[8]: fam2label = build_labels(train_targets)
    There are 17930 labels.
```

Data exploration and analysis: a few plots to understand the data at hand



One can see that the provided dataset is **heavily imbalanced**: some families have up to roughly 3k samples while others only have around ten samples.

```
[10]: # Plot the distribution of sequences' lengths

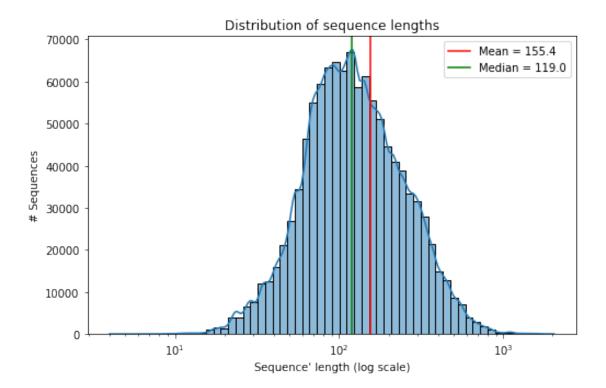
f, ax = plt.subplots(figsize=(8, 5))

sequence_lengths = train_data.str.len()
median = sequence_lengths.median()
mean = sequence_lengths.mean()

sns.histplot(sequence_lengths.values, kde=True, log_scale=True, bins=60, ax=ax)

ax.axvline(mean, color='r', linestyle='-', label=f"Mean = {mean:.1f}")
ax.axvline(median, color='g', linestyle='-', label=f"Median = {median:.1f}")

plt.title("Distribution of sequence lengths")
plt.xlabel("Sequence' length (log scale)")
plt.ylabel("# Sequences")
plt.legend(loc="best")
plt.show()
```



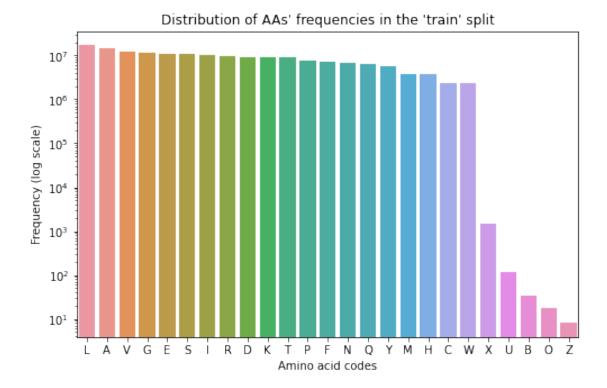
One can observe that the sequences' lengths are quite different over the dataset.

```
[11]: def get_amino_acid_frequencies(data):
    aa_counter = Counter()

    for sequence in data:
        aa_counter.update(sequence)

    return pd.DataFrame({'AA': list(aa_counter.keys()), 'Frequency':
        →list(aa_counter.values())})
```





Eventually, one can observe that some amino acids (X, U, B, O, Z) are quite rare compared to other ones. We choose to consider those rare amino acids as unknown (<unk>) amino acids but this can be changed.

```
[13]: def build_vocab(data):
    # Build the vocabulary
    voc = set()
    rare_AAs = {'X', 'U', 'B', 'O', 'Z'}
    for sequence in data:
        voc.update(sequence)

    unique_AAs = sorted(voc - rare_AAs)

# Build the mapping
    word2id = {w: i for i, w in enumerate(unique_AAs, start=2)}
    word2id['<pad>'] = 0
    word2id['<unk>'] = 1

    return word2id
```

```
[14]: word2id = build_vocab(train_data)
[15]: word2id
[15]: {'A': 2,
       'C': 3,
       'D': 4,
       'E': 5,
       'F': 6,
       'G': 7,
       'H': 8,
       'I': 9,
       'K': 10,
       'L': 11,
       'M': 12,
       'N': 13,
       'P': 14,
       'Q': 15,
       'R': 16,
       'S': 17,
       'T': 18,
       'V': 19,
       'W': 20,
       'Y': 21,
       '<pad>': 0,
       '<unk>': 1}
[16]: print(f"AA dictionary formed. The length of dictionary is: {len(word2id)}.")
     AA dictionary formed. The length of dictionary is: 22.
[17]: class SequenceDataset(torch.utils.data.Dataset):
          def __init__(self, word2id, fam2label, max_len, data_path, split):
              self.word2id = word2id
              self.fam2label = fam2label
              self.max_len = max_len
              self.data, self.label = reader(split, data_path)
          def __len__(self):
              return len(self.data)
          def __getitem__(self, index):
              seq = self.preprocess(self.data.iloc[index])
              label = self.fam2label.get(self.label.iloc[index], self.
       →fam2label['<unk>'])
```

```
return {'sequence': seq, 'target' : label}
          def preprocess(self, text):
              seq = []
              # Encode into IDs
              for word in text[:self.max_len]:
                  seq.append(self.word2id.get(word, self.word2id['<unk>']))
              # Pad to maximal length
              if len(seq) < self.max_len:</pre>
                  seq += [self.word2id['<pad>'] for _ in range(self.max_len -__
       \rightarrowlen(seq))]
              # Convert list into tensor
              seq = torch.from_numpy(np.array(seq))
              # One-hot encode
              one_hot_seq = torch.nn.functional.one_hot(seq, num_classes=len(self.
       →word2id), )
              # Permute channel (one-hot) dim first
              one_hot_seq = one_hot_seq.permute(1,0)
              return one_hot_seq
[18]: seq_max_len = 120
[19]: train_dataset = SequenceDataset(word2id, fam2label, seq_max_len, data_dir,__
      dev_dataset = SequenceDataset(word2id, fam2label, seq_max_len, data_dir, "dev")
      test_dataset = SequenceDataset(word2id, fam2label, seq_max_len, data_dir,__
       →"test")
[20]: # Health check: test the dataset
      next(iter(train_dataset))['sequence'].shape
[20]: torch.Size([22, 120])
[21]: batch_size = 1
[22]: num_workers = 0
      dataloaders = {}
      dataloaders['train'] = torch.utils.data.DataLoader(
          train_dataset,
```

```
batch_size=batch_size,
          shuffle=True,
          num_workers=num_workers,
      dataloaders['dev'] = torch.utils.data.DataLoader(
          dev_dataset,
          batch_size=batch_size,
          shuffle=False,
          num_workers=num_workers,
      )
      dataloaders['test'] = torch.utils.data.DataLoader(
          test_dataset,
          batch_size=batch_size,
          shuffle=False,
          num_workers=num_workers,
[23]: # Health check with the dataloader
      batch = next(iter(dataloaders['test']))
      batch['sequence'].shape, batch['target'].shape
[23]: (torch.Size([1, 22, 120]), torch.Size([1]))
     0.1.2 Step 2: Build the classification model
[24]: num_classes = len(fam2label)
[25]: class Lambda(torch.nn.Module):
          def __init__(self, func):
              super().__init__()
              self.func = func
          def forward(self, x):
              return self.func(x)
      class ResidualBlock(torch.nn.Module):
          The residual block used by ProtCNN (https://www.biorxiv.org/content/10.1101/
       \hookrightarrow 626507v3. full).
          Args:
              in_channels: The number of channels (feature maps) of the incoming_
       \hookrightarrow embedding
               out_channels: The number of channels after the first convolution
               dilation: Dilation rate of the first convolution
```

```
def __init__(self, in_channels, out_channels, dilation=1):
       super().__init__()
       # Initialize the required layers
       self.skip = torch.nn.Sequential()
       self.bn1 = torch.nn.BatchNorm1d(in_channels)
       self.conv1 = torch.nn.Conv1d(in_channels=in_channels,_
→out_channels=out_channels,
                              kernel_size=3, bias=False, dilation=dilation,__
→padding=dilation)
       self.bn2 = torch.nn.BatchNorm1d(out_channels)
       self.conv2 = torch.nn.Conv1d(in_channels=out_channels,__
→out_channels=out_channels,
                              kernel_size=3, bias=False, padding=1)
   def forward(self, x):
       # Execute the required layers and functions
       activation = F.relu(self.bn1(x))
       x1 = self.conv1(activation)
       x2 = self.conv2(F.relu(self.bn2(x1)))
       return x2 + self.skip(x)
```

```
[26]: class ProtCNN(pl.LightningModule):
          def __init__(self, num_classes):
              super().__init__()
              self.model = torch.nn.Sequential(
                  torch.nn.Conv1d(22, 128, kernel_size=1, padding=0, bias=False),
                  ResidualBlock(128, 128, dilation=2),
                  ResidualBlock(128, 128, dilation=3),
                  torch.nn.MaxPool1d(3, stride=2, padding=1),
                  Lambda(lambda x: x.flatten(start_dim=1)),
                  torch.nn.Linear(7680, num_classes)
              )
              self.train_acc = torchmetrics.Accuracy()
              self.valid_acc = torchmetrics.Accuracy()
          def forward(self, x):
              return self.model(x.float())
          def training_step(self, batch, batch_idx):
              x, y = batch['sequence'], batch['target']
              y_hat = self(x)
```

```
loss = F.cross_entropy(y_hat, y)
              self.log('train_loss', loss, on_step=True, on_epoch=True)
              pred = torch.argmax(y_hat, dim=1)
              self.train_acc(pred, y)
              self.log('train_acc', self.train_acc, on_step=True, on_epoch=True)
              return loss
          def validation_step(self, batch, batch_idx):
              x, y = batch['sequence'], batch['target']
              y_hat = self(x)
              pred = torch.argmax(y_hat, dim=1)
              acc = self.valid_acc(pred, y)
              self.log('valid_acc', self.valid_acc, on_step=False, on_epoch=True)
              return acc
          def configure_optimizers(self):
              optimizer = torch.optim.SGD(self.parameters(), lr=1e-2, momentum=0.9, u
       →weight_decay=1e-2)
              lr scheduler = torch.optim.lr scheduler.MultiStepLR(optimizer,...)
       →milestones=[5, 8, 10, 12, 14, 16, 18, 20], gamma=0.9)
              return {
                  "optimizer": optimizer,
                  "lr scheduler": lr scheduler,
              }
[27]: prot_cnn = ProtCNN(num_classes)
[28]: prot_cnn
[28]: ProtCNN(
        (model): Sequential(
          (0): Conv1d(22, 128, kernel_size=(1,), stride=(1,), bias=False)
          (1): ResidualBlock(
            (skip): Sequential()
            (bn1): BatchNorm1d(128, eps=1e-05, momentum=0.1, affine=True,
      track_running_stats=True)
            (conv1): Conv1d(128, 128, kernel_size=(3,), stride=(1,), padding=(2,),
      dilation=(2,), bias=False)
            (bn2): BatchNorm1d(128, eps=1e-05, momentum=0.1, affine=True,
      track_running_stats=True)
            (conv2): Conv1d(128, 128, kernel_size=(3,), stride=(1,), padding=(1,),
      bias=False)
          )
```

```
(2): ResidualBlock(
            (skip): Sequential()
            (bn1): BatchNorm1d(128, eps=1e-05, momentum=0.1, affine=True,
      track_running_stats=True)
            (conv1): Conv1d(128, 128, kernel_size=(3,), stride=(1,), padding=(3,),
      dilation=(3,), bias=False)
            (bn2): BatchNorm1d(128, eps=1e-05, momentum=0.1, affine=True,
      track_running_stats=True)
            (conv2): Conv1d(128, 128, kernel_size=(3,), stride=(1,), padding=(1,),
     bias=False)
          (3): MaxPool1d(kernel_size=3, stride=2, padding=1, dilation=1,
      ceil mode=False)
          (4): Lambda()
          (5): Linear(in_features=7680, out_features=17930, bias=True)
        (train_acc): Accuracy()
        (valid_acc): Accuracy()
      )
[29]: pl.seed_everything(0)
     Global seed set to 0
[29]: 0
[30]: # Health check: test that the network works on a single mini-batch
      batch = next(iter(dataloaders['train']))
      print(batch["sequence"].shape)
      prot_cnn(batch["sequence"]).shape
     torch.Size([1, 22, 120])
[30]: torch.Size([1, 17930])
[31]: gpus = 0
      epochs = 25
[32]: trainer = pl.Trainer(gpus=gpus, max_epochs=epochs)
     GPU available: False, used: False
     TPU available: False, using: 0 TPU cores
     IPU available: False, using: 0 IPUs
[33]: trainer.fit(prot_cnn, dataloaders['train'], dataloaders['dev'])
       Name
                   | Type
                                | Params
```

137 M Trainable params
0 Non-trainable params

137 M Total params

551.683 Total estimated model params size (MB)

Validation sanity check: 0it [00:00, ?it/s]

/Users/theomeborckinstadeep/Desktop/experiments/venv/lib/python3.7/site-packages/pytorch_lightning/trainer/data_loading.py:112: UserWarning: The dataloader, val_dataloader 0, does not have many workers which may be a bottleneck. Consider increasing the value of the `num_workers` argument` (try 8 which is the number of cpus on this machine) in the `DataLoader` init to improve performance.

f"The dataloader, {name}, does not have many workers which may be a bottleneck."

Global seed set to 0

/Users/theomeborckinstadeep/Desktop/experiments/venv/lib/python3.7/site-packages/pytorch_lightning/trainer/data_loading.py:112: UserWarning: The dataloader, train_dataloader, does not have many workers which may be a bottleneck. Consider increasing the value of the `num_workers` argument` (try 8 which is the number of cpus on this machine) in the `DataLoader` init to improve performance.

f"The dataloader, {name}, does not have many workers which may be a bottleneck."

Training: Oit [00:00, ?it/s]

/Users/theomeborckinstadeep/Desktop/experiments/venv/lib/python3.7/site-packages/pytorch_lightning/trainer/trainer.py:685: UserWarning: Detected KeyboardInterrupt, attempting graceful shutdown...

rank_zero_warn("Detected KeyboardInterrupt, attempting graceful shutdown...")

0.1.3 Step 3: Visualize the training

```
[29]: # Load the TensorBoard notebook extension %load_ext tensorboard
```

```
[30]: %tensorboard --logdir "./lightning_logs/"
```

<IPython.core.display.HTML object>