

Trainer

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<u>Trainer</u> is a complete training and evaluation loop for Transformers' PyTorch models. Plug a model, preprocessor, dataset, and training arguments into <u>Trainer</u> and let it handle the rest to start training faster.

<u>Trainer</u> is also powered by <u>Accelerate</u>, a library for handling large models for distributed training.

This guide will show you how <u>Trainer</u> works and how to customize it for your use case with a callback.

!pip install accelerate --upgrade

<u>Trainer</u> contains all the necessary components of a training loop.

- 1. calculate the loss from a training step
- 2. calculate the gradients with the <u>backward</u> method
- 3. update the weights based on the gradients
- 4. repeat until the predetermined number of epochs is reached

Manually coding this training loop everytime can be inconvenient or a barrier if you're just getting started with machine learning. <u>Trainer</u> abstracts this process, allowing you to focus on the model, dataset, and training design choices.

Configure your training with hyperparameters and options from <u>TrainingArguments</u> which supports many features such as distributed training, torch.compile, mixed precision training, and saving the model to the Hub.

The number of available parameters available in <u>TrainingArguments</u> may be intimidating at first. If there is a specific hyperparameter or feature you want to use, try searching for it directly.

Otherwise, feel free to start with the default values and gradually customize them as you become more familiar with the training process.

The example below demonstrates an example of <u>TrainingArguments</u> that evaluates and saves the model at the end of each epoch. It also loads the best model found during training and pushes it to the Hub.

```
from transformers import TrainingArguments

training_args = TrainingArguments(
    output_dir="your-model",
    learning_rate=2e-5,
    per_device_train_batch_size=16,
    per_device_eval_batch_size=16,
    num_train_epochs=2,
    weight_decay=0.01,
    eval_strategy="epoch",
    save_strategy="epoch",
    load_best_model_at_end=True,
    push_to_hub=True,
)
```

Pass your model, dataset, preprocessor, and <u>TrainingArguments</u> to <u>Trainer</u>, and call <u>train()</u> to start training.

Refer to the Fine-tuning guide for a more complete overview of the training process.

```
from transformers import Trainer

trainer = Trainer(
    model=model,
    args=training_args,
    train_dataset=dataset["train"],
    eval_dataset=dataset["test"],
    processing_class=tokenizer,
    data_collator=data_collator,
    compute_metrics=compute_metrics,
)

trainer.train()
```

Checkpoints

<u>Trainer</u> saves checkpoints (the optimizer state is not saved by default) to the directory in output_dir in <u>TrainingArguments</u> to a subfolder named checkpoint-000. The number at the end is the training step at which the checkpoint was saved.

Saving checkpoints are useful for resuming training or recovering your training progress if you encounter an error. Set the resume_from_checkpoint parameter in train() to resume training from the last checkpoint or a specific checkpoint.

latest checkpoint (specific checkpoint

trainer.train(resume_from_checkpoint=True)

Checkpoints can be saved to the Hub by setting push_to_hub=True in <u>TrainingArguments</u>. The default method ("every_save") saves a checkpoint to the Hub every time a model is saved, which is typically the final model at the end of training. Some other options for deciding how to save checkpoints to the Hub include the following.

- hub_strategy="end" only pushes a checkpoint when save model() is called
- hub_strategy="checkpoint" pushes the latest checkpoint to a subfolder named lastcheckpoint from which training can be resumed
- hub_strategy="all_checkpoints" pushes all checkpoints to the Hub with one checkpoint per subfolder in your model repository

<u>Trainer</u> attempts to maintain the same Python, NumPy, and PyTorch RNG states when you resume training from a checkpoint. But PyTorch has various non-deterministic settings which can't guarantee the RNG states are identical. To enable full determinism, refer to the <u>Controlling sources of randomness</u> guide to learn what settings to adjust to make training fully deterministic (some settings may result in slower training).

Logging

<u>Trainer</u> is set to logging. INFO by default to report errors, warnings, and other basic information. Use log_level() to change the logging level and log verbosity.

The example below sets the main code and modules to use the same log level.

```
logger = logging.getLogger(__name__)

logging.basicConfig(
    format="%(asctime)s - %(levelname)s - %(name)s - %(message)s",
    datefmt="%m/%d/%Y %H:%M:%S",
    handlers=[logging.StreamHandler(sys.stdout)],
)

log_level = training_args.get_process_log_level()
logger.setLevel(log_level)
datasets.utils.logging.set_verbosity(log_level)
transformers.utils.logging.set_verbosity(log_level)
trainer = Trainer(...)
```

In a distributed environment, <u>Trainer</u> replicas are set to logging.WARNING to only report errors and warnings. Use log_level_replica() to change the logging level and log verbosity. To configure the log level for each node, use log_on_each_node() to determine whether to use a specific log level on each node or only the main node.

Use different combinations of log_level and log_level_replica to configure what gets logged on each node.

```
single node multi-node

my_app.py ... --log_level warning --log_level_replica error
```

The log level is separately set for each node in the <code>__init__()</code> method. Consider setting this sooner if you're using other Transformers functionalities before creating the <u>Trainer</u> instance.

Customize

Tailor <u>Trainer</u> to your use case by subclassing or overriding its methods to support the functionality you want to add or use, without rewriting the entire training loop from scratch. The table below lists some of the methods that can be customized.

method	description
get train dataloader()	create a training DataLoader
get eval dataloader()	create an evaluation DataLoader
get test dataloader()	create a test DataLoader
<u>log()</u>	log information about the training process
create optimizer and scheduler()	create an optimizer and learning rate scheduler (can also be separately customized with create_optimizer() and create_scheduler() if they weren't passed ininit)
compute loss()	compute the loss of a batch of training inputs
training step()	perform the training step
<u>prediction_step()</u>	perform the prediction and test step
<u>evaluate()</u>	evaluate the model and return the evaluation metric
<u>predict()</u>	make a prediction (with metrics if labels are available) on the test set

For example, to use weighted loss, rewrite <u>compute loss()</u> inside <u>Trainer</u>.

```
from torch import nn
from transformers import Trainer

class CustomTrainer(Trainer):
    def compute_loss(self, model: nn.Module, inputs: dict[str, Union[torch.Tensor, Any
        labels = inputs.pop("labels")
        # forward pass
        outputs = model(**inputs)
        logits = outputs.get("logits")
        # compute custom loss for 3 labels with different weights
        reduction = "sum" if num_items_in_batch is not None else "mean"
        loss_fct = nn.CrossEntropyLoss(weight=torch.tensor([1.0, 2.0, 3.0], device=mod
        loss = loss_fct(logits.view(-1, self.model.config.num_labels), labels.view(-1)
        if num_items_in_batch is not None:
            loss = loss / num_items_in_batch
            return (loss, outputs) if return_outputs else loss
```

Callbacks

<u>Callbacks</u> are another way to customize <u>Trainer</u>, but they don't change anything *inside the training loop*. Instead, a callback inspects the training loop state and executes some action (early stopping, logging, etc.) depending on the state. For example, you can't implement a custom loss function with a callback because that requires overriding compute loss().

To use a callback, create a class that inherits from <u>TrainerCallback</u> and implements the functionality you want. Then pass the callback to the callback parameter in <u>Trainer</u>. The example below implements an early stopping callback that stops training after 10 steps.

```
from transformers import TrainerCallback, Trainer
class EarlyStoppingCallback(TrainerCallback):
    def __init__(self, num_steps=10):
        self.num_steps = num_steps
    def on_step_end(self, args, state, control, **kwargs):
        if state.global_step >= self.num_steps:
            return {"should training stop": True}
        else:
            return {}
trainer = Trainer(
    model=model,
    args=training_args,
    train_dataset=dataset["train"],
    eval_dataset=dataset["test"],
    processing_class=tokenizer,
    data_collator=data_collator,
    compute_metrics=compute_metrics,
    callbacks=[EarlyStoppingCallback()],
)
```

Accelerate

<u>Accelerate</u> is a library that simplifies training in distributed environments and across different hardware. Its integration with <u>Trainer</u> means <u>Trainer</u> supports distributed training frameworks like <u>Fully Sharded Data Parallel (FSDP)</u> and <u>DeepSpeed</u>.

Learn more about FSDP sharding strategies, CPU offloading, and more with <u>Trainer</u> in the <u>Fully</u>

<u>Sharded Data Parallel</u> guide.

To use Accelerate with <u>Trainer</u>, run the <u>accelerate config</u> command to configure your training environment. This command creates a config_file.yaml file that stores the configuration settings of your training environment and it's used whenever you launch your training script. Some example distributed training configurations are shown below.

```
DeepSpeed with Accelerate
                      FullyShardedDataParallel
DistributedDataParallel
                                              DeepSpeed
                                                          plugin
compute_environment: LOCAL_MACHINE
distributed type: MULTI GPU
downcast_bf16: 'no'
gpu_ids: all
machine_rank: 0 #change rank as per the node
main_process_ip: 192.168.20.1
main_process_port: 9898
main_training_function: main
mixed_precision: fp16
num_machines: 2
num processes: 8
rdzv_backend: static
same network: true
tpu env: []
tpu_use_cluster: false
tpu_use_sudo: false
use_cpu: false
```

Run <u>accelerate launch</u> to start training with the configurations set in config_file.yaml. This file is saved to the Accelerate cache folder and automatically loaded when you run accelerate_launch.

The example below launches the <u>run_glue.py</u> script with the FSDP configuration shown earlier. Parameters from the config_file.yaml file can also be directly set in the command line.

```
accelerate launch \
    ./examples/pytorch/text-classification/run_glue.py \
    --model_name_or_path google-bert/bert-base-cased \
    --task_name $TASK_NAME \
    --do_train \
    --do_eval \
    --max_seq_length 128 \
    --per_device_train_batch_size 16 \
```

```
--learning_rate 5e-5 \
--num_train_epochs 3 \
--output_dir /tmp/$TASK_NAME/ \
--overwrite_output_dir
```

Refer to the <u>Launching your Accelerate scripts</u> tutorial to learn more about accelerate_launch and custom configurations.

Optimizations

<u>Trainer</u> supports various optimizations to improve *training* performance - reduce memory and increase training speed - and *model* performance.

torch.compile

<u>torch.compile</u> can significantly speed up training and reduce computational overhead.

Configure your torch.compile settings in <u>TrainingArguments</u>. Set torch_compile to True, and select a backend and compile mode.

```
from transformers import TrainingArguments

training_args = TrainingArguments(
    torch_compile=True,
    torch_compile_backend="inductor",
    torch_compile_mode="default",
    ...,
)
```

GaLore

<u>Gradient Low-Rank Projection (GaLore)</u> significantly reduces memory usage when training large language models (LLMs). One of GaLores key benefits is *full-parameter* learning, unlike low-rank adaptation methods like <u>LoRA</u>, which produces better model performance.

Install the GaLore and TRL libraries.

```
pip install galore-torch trl
```

Pick a GaLore optimizer ("galore_adamw", "galore_adafactor", "galore_adamw_8bit") and pass it to the optim parameter in <u>trl.SFTConfig</u>. Use the optim_target_modules parameter to specify which modules to adapt (can be a list of strings, regex, or a full path).

Extra parameters supported by GaLore, rank, update_proj_gap, and scale, should be passed to the optim_args parameter in trl.SFTConfig.

The example below enables GaLore with <u>SFTTrainer</u> that targets the attn and mlp layers with regex.

It can take some time before training starts (~3 minutes for a 2B model on a NVIDIA A100).

(GaLore optimizer) GaLore optimizer with layerwise optimization

```
import datasets
from trl import SFTConfig, SFTTrainer
train_dataset = datasets.load_dataset('imdb', split='train')
args = SFTConfig(
   output_dir="./test-galore",
   max_steps=100,
    optim="galore_adamw",
    optim_target_modules=[r".*.attn.*", r".*.mlp.*"],
    optim_args="rank=64, update_proj_gap=100, scale=0.10",
    gradient_checkpointing=True,
)
trainer = SFTTrainer(
   model="google/gemma-2b",
    args=args,
    train_dataset=train_dataset,
)
trainer.train()
```

Only linear layers that are considered GaLore layers can be trained with low-rank decomposition. The rest of the model layers are optimized in the usual way.

Liger

<u>Liger Kernel</u> is a collection of layers such as RMSNorm, RoPE, SwiGLU, CrossEntropy, FusedLinearCrossEntropy, and more that have been fused into a single Triton kernel for training

LLMs. These kernels are also compatible with FlashAttention, FSDP, and DeepSpeed. As a result, Liger Kernel can increase multi-GPU training throughput and reduce memory usage. This is useful for multi-head training and supporting larger vocabulary sizes, larger batch sizes, and longer context lengths.

```
pip install liger-kernel
```

Enable Liger Kernel for training by setting use_liger_kernel=True in <u>TrainingArguments</u>. This patches the corresponding layers in the model with Ligers kernels.

Liger Kernel supports Llama, Gemma, Mistral, and Mixtral models. Refer to the <u>patching</u> list for the latest list of supported models.

```
from transformers import TrainingArguments

training_args = TrainingArguments(
    output_dir="your-model",
    learning_rate=2e-5,
    per_device_train_batch_size=16,
    per_device_eval_batch_size=16,
    num_train_epochs=2,
    weight_decay=0.01,
    eval_strategy="epoch",
    save_strategy="epoch",
    load_best_model_at_end=True,
    push_to_hub=True,
    use_liger_kernel=True
)
```

You can also configure which specific kernels to apply using the liger_kernel_config parameter. This dict is passed as keyword arguments to the _apply_liger_kernel_to_instance function, allowing fine-grained control over kernel usage. Available options vary by model but typically include: rope, swiglu, cross_entropy, fused_linear_cross_entropy, rms_norm, etc.

```
from transformers import TrainingArguments
# Apply only specific kernels
```

```
training_args = TrainingArguments(
    output_dir="your-model",
    learning_rate=2e-5,
    per device train batch size=16,
    per_device_eval_batch_size=16,
    num train epochs=2,
    weight_decay=0.01,
    eval strategy="epoch",
    save strategy="epoch",
    load_best_model_at_end=True,
    push_to_hub=True,
    use_liger_kernel=True,
    liger_kernel_config={
        "rope": True,
        "cross_entropy": True,
        "rms_norm": False, # Don't apply Liger's RMSNorm kernel
        "swiglu": True,
    3
)
```

NEFTune

<u>NEFTune</u> adds noise to the embedding vectors during training to improve model performance. Enable it in <u>Trainer</u> with the neftune_noise_alpha parameter in <u>TrainingArguments</u> to control how much noise is added.

```
from transformers import TrainingArguments, Trainer

training_args = TrainingArguments(..., neftune_noise_alpha=0.1)
trainer = Trainer(..., args=training_args)
```

The original embedding layer is restored after training to avoid any unexpected behavior.

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← Inference server backends

Fine-tuning →