Depending on the dataset and your use case, your test dataset may contain labels. In that case, this method will also return metrics, like in evaluate().

If your predictions or labels have different sequence lengths (for instance because you're doing dynamic padding in a token classification task) the predictions will be padded (on the right) to allow for concatenation into one array. The padding index is -100.

Returns: NamedTuple A namedtuple with the following keys:

- •predictions (np.ndarray): The predictions on test\_dataset.
- •label\_ids (np.ndarray, optional): The labels (if the dataset contained some).
- metrics (Dict[str, float], optional): The potential dictionary of metrics (if the dataset contained labels).

## **TrainingArguments**

## class transformers.TrainingArguments

( output dir: str, overwrite output dir: bool = False, do train: bool = False, do eval: bool = False, do\_predict: bool = False, eval\_strategy: Union = 'no', prediction\_loss\_only: bool = False, per\_device\_train\_batch\_size: int = 8, per\_device\_eval\_batch\_size: int = 8, per\_gpu\_train\_batch\_size: Optional = None, per\_gpu\_eval\_batch\_size: Optional = None, gradient\_accumulation\_steps: int = 1, eval\_accumulation\_steps: Optional = None, eval\_delay: Optional = 0, learning\_rate: float = 5e-05, weight\_decay: float = 0.0, adam\_beta1: float = 0.9, adam\_beta2: float = 0.999, adam\_epsilon: float = 1e-08, max\_grad\_norm: float = 1.0, num\_train\_epochs: float = 3.0, max\_steps: int = -1, lr\_scheduler\_type: Union = 'linear', lr\_scheduler\_kwargs: Union = <factory>, warmup\_ratio: float = 0.0, warmup\_steps: int = 0, log\_level: Optional = 'passive', log\_level\_replica: Optional = 'warning', log\_on\_each\_node: bool = True, logging\_dir: Optional = None, logging\_strategy: Union = 'steps', logging\_first\_step: bool = False, logging\_steps: float = 500, logging\_nan\_inf\_filter: bool = True, save\_strategy: Union = 'steps', save\_steps: float = 500, save\_total\_limit: Optional = None, save\_safetensors: Optional = True, save\_on\_each\_node: bool = False, save\_only\_model: bool = False, restore\_callback\_states\_from\_checkpoint: bool = False, no\_cuda: bool = False, use\_cpu: bool = False, use\_mps\_device: bool = False, seed: int = 42, data\_seed: Optional = None, jit\_mode\_eval: bool = False, use\_ipex: bool = False, bf16: bool = False, fp16: bool = False, fp16\_opt\_level: str = '01', half\_precision\_backend: str = 'auto', bf16\_full\_eval: bool = False, fp16\_full\_eval: bool = False, tf32: Optional = None, local\_rank: int = -1, ddp\_backend: Optional = None, tpu\_num\_cores: Optional = None, tpu\_metrics\_debug: bool = False, debug: Union = '', dataloader\_drop\_last: bool = False, eval\_steps: Optional = None, dataloader\_num\_workers: int = 0, dataloader\_prefetch\_factor: Optional = None, past\_index: int = -1, run\_name: Optional = None, disable\_tqdm: Optional = None, remove\_unused\_columns: Optional = True, label\_names: Optional = None, load\_best\_model\_at\_end: Optional = False, metric\_for\_best\_model: Optional = None, greater\_is\_better: Optional = None, ignore\_data\_skip: bool = False, fsdp: Union = '', fsdp\_min\_num\_params: int = 0, fsdp\_config: Union = None, fsdp\_transformer\_layer\_cls\_to\_wrap: Optional = None, accelerator\_config: Union = None, deepspeed: Union = None, label\_smoothing\_factor: float = 0.0, optim: Union = 'adamw torch', optim\_args: Optional = None, adafactor: bool = False, group\_by\_length: bool = False, length\_column\_name: Optional = 'length', report\_to: Union = None, ddp\_find\_unused\_parameters:

```
Optional = None, ddp_bucket_cap_mb: Optional = None, ddp_broadcast_buffers: Optional = None, dataloader_pin_memory: bool = True, dataloader_persistent_workers: bool = False, skip_memory_metrics: bool = True, use_legacy_prediction_loop: bool = False, push_to_hub: bool = False, resume_from_checkpoint: Optional = None, hub_model_id: Optional = None, hub_strategy: Union = 'every_save', hub_token: Optional = None, hub_private_repo: bool = False, hub_always_push: bool = False, gradient_checkpointing: bool = False, gradient_checkpointing_kwargs: Union = None, include_inputs_for_metrics: bool = False, eval_do_concat_batches: bool = True, fp16_backend: str = 'auto', evaluation_strategy: Union = None, push_to_hub_model_id: Optional = None, push_to_hub_organization: Optional = None, push_to_hub_token: Optional = None, mp_parameters: str = '', auto_find_batch_size: bool = False, full_determinism: bool = False, torchdynamo: Optional = None, ray_scope: Optional = 'last', ddp_timeout: Optional = 1800, torch_compile: bool = False, torch_compile_backend: Optional = None, torch_compile_mode: Optional = None, dispatch_batches: Optional = None, split_batches: Optional = None, include_tokens_per_second: Optional = False, include_num_input_tokens_seen: Optional = False, neftune_noise_alpha: Optional = None, optim_target_modules: Union = None, batch_eval_metrics: bool = False)
```

## **Parameters**

- output\_dir (str) The output directory where the model predictions and checkpoints will be written.
- **overwrite\_output\_dir** (bool, *optional*, defaults to False) If True, overwrite the content of the output directory. Use this to continue training if output\_dir points to a checkpoint directory.
- **do\_train** (bool, *optional*, defaults to False) Whether to run training or not. This argument is not directly used by <u>Trainer</u>, it's intended to be used by your training/evaluation scripts instead. See the <u>example scripts</u> for more details.
- **do\_eval** (boo1, *optional*) Whether to run evaluation on the validation set or not. Will be set to True if eval\_strategy is different from "no". This argument is not directly used by <u>Trainer</u>, it's intended to be used by your training/evaluation scripts instead. See the <u>example scripts</u> for more details.
- do\_predict (bool, optional, defaults to False) Whether to run predictions on the test set or not.
   This argument is not directly used by <u>Trainer</u>, it's intended to be used by your training/evaluation scripts instead. See the <u>example scripts</u> for more details.
- **eval\_strategy** (str or <u>IntervalStrategy</u>, *optional*, defaults to "no") The evaluation strategy to adopt during training. Possible values are:
  - "no": No evaluation is done during training.
  - "steps": Evaluation is done (and logged) every eval\_steps.
  - "epoch": Evaluation is done at the end of each epoch.
- **prediction\_loss\_only** (boo1, *optional*, defaults to False) When performing evaluation and generating predictions, only returns the loss.
- per\_device\_train\_batch\_size (int, optional, defaults to 8) The batch size per GPU/XPU/TPU/MPS/NPU core/CPU for training.
- per\_device\_eval\_batch\_size (int, optional, defaults to 8) The batch size per GPU/XPU/TPU/MPS/NPU core/CPU for evaluation.

**gradient\_accumulation\_steps** (int, *optional*, defaults to 1) — Number of updates steps to accumulate the gradients for, before performing a backward/update pass.

When using gradient accumulation, one step is counted as one step with backward pass. Therefore, logging, evaluation, save will be conducted every gradient\_accumulation\_steps \* xxx\_step training examples.

- eval\_accumulation\_steps (int, optional) Number of predictions steps to accumulate the output tensors for, before moving the results to the CPU. If left unset, the whole predictions are accumulated on GPU/NPU/TPU before being moved to the CPU (faster but requires more memory).
- eval\_delay (float, optional) Number of epochs or steps to wait for before the first evaluation can be performed, depending on the eval\_strategy.
- **learning\_rate** (float, optional, defaults to 5e-5) The initial learning rate for <u>AdamW</u> optimizer.
- weight\_decay (float, optional, defaults to 0) The weight decay to apply (if not zero) to all layers except all bias and LayerNorm weights in <u>AdamW</u> optimizer.
- adam\_beta1 (float, optional, defaults to 0.9) The beta1 hyperparameter for the <u>AdamW</u> optimizer.
- adam\_beta2 (float, optional, defaults to 0.999) The beta2 hyperparameter for the <u>AdamW</u> optimizer.
- adam\_epsilon (float, optional, defaults to 1e-8) The epsilon hyperparameter for the <u>AdamW</u> optimizer.
- max\_grad\_norm (float, optional, defaults to 1.0) Maximum gradient norm (for gradient clipping).
- **num\_train\_epochs**(float, *optional*, defaults to 3.0) Total number of training epochs to perform (if not an integer, will perform the decimal part percents of the last epoch before stopping training).
- max\_steps (int, optional, defaults to -1) If set to a positive number, the total number of training steps to perform. Overrides num\_train\_epochs. For a finite dataset, training is reiterated through the dataset (if all data is exhausted) until max\_steps is reached.
- **lr\_scheduler\_type** (str or <u>SchedulerType</u>, *optional*, defaults to "linear") The scheduler type to use. See the documentation of <u>SchedulerType</u> for all possible values.
- **lr\_scheduler\_kwargs** ('dict', *optional*, defaults to {}) The extra arguments for the lr\_scheduler. See the documentation of each scheduler for possible values.
- warmup\_ratio (float, optional, defaults to 0.0) Ratio of total training steps used for a linear warmup from 0 to learning\_rate.
- warmup\_steps (int, optional, defaults to 0) Number of steps used for a linear warmup from 0 to learning\_rate. Overrides any effect of warmup\_ratio.
- **log\_level** (str, *optional*, defaults to passive) Logger log level to use on the main process.

  Possible choices are the log levels as strings: 'debug', 'info', 'warning', 'error' and 'critical', plus a 'passive' level which doesn't set anything and keeps the current log level for the Transformers library (which will be "warning" by default).

• log\_level\_replica (str, optional, defaults to "warning") — Logger log level to use on replicas. Same choices as log\_level"

- **log\_on\_each\_node** (bool, *optional*, defaults to True) In multinode distributed training, whether to log using log\_level once per node, or only on the main node.
- logging\_dir (str, optional) <u>TensorBoard</u> log directory. Will default to
   \*output dir/runs/CURRENT\_DATETIME\_HOSTNAME\*.
- **logging\_strategy** (str or <u>IntervalStrategy</u>, *optional*, defaults to "steps") The logging strategy to adopt during training. Possible values are:
  - "no": No logging is done during training.
  - "epoch": Logging is done at the end of each epoch.
  - "steps": Logging is done every logging\_steps.
- logging\_first\_step (bool, optional, defaults to False) Whether to log the first global\_step or not.
- **logging\_steps** (int or float, *optional*, defaults to 500) Number of update steps between two logs if logging\_strategy="steps". Should be an integer or a float in range [0,1). If smaller than 1, will be interpreted as ratio of total training steps.
- **logging\_nan\_inf\_filter** (bool, *optional*, defaults to True) Whether to filter nan and inf losses for logging. If set to True the loss of every step that is nan or inf is filtered and the average loss of the current logging window is taken instead.

logging\_nan\_inf\_filter only influences the logging of loss values, it does not change the behavior the gradient is computed or applied to the model.

- save\_strategy (str or <u>IntervalStrategy</u>, optional, defaults to "steps") The checkpoint save strategy to adopt during training. Possible values are:
  - "no": No save is done during training.
  - "epoch": Save is done at the end of each epoch.
  - "steps": Save is done every save\_steps.
- save\_steps (int or float, optional, defaults to 500) Number of updates steps before two checkpoint saves if save\_strategy="steps". Should be an integer or a float in range [0,1). If smaller than 1, will be interpreted as ratio of total training steps.
- save\_total\_limit (int, optional) If a value is passed, will limit the total amount of checkpoints. Deletes the older checkpoints in output\_dir. When load\_best\_model\_at\_end is enabled, the "best" checkpoint according to metric\_for\_best\_model will always be retained in addition to the most recent ones. For example, for save\_total\_limit=5 and load\_best\_model\_at\_end, the four last checkpoints will always be retained alongside the best model. When save\_total\_limit=1 and load\_best\_model\_at\_end, it is possible that two checkpoints are saved: the last one and the best one (if they are different).

• save\_safetensors (bool, optional, defaults to True) — Use <u>safetensors</u> saving and loading for state dicts instead of default torch.load and torch.save.

- **save\_on\_each\_node** (bool, *optional*, defaults to False) When doing multi-node distributed training, whether to save models and checkpoints on each node, or only on the main one.
  - This should not be activated when the different nodes use the same storage as the files will be saved with the same names for each node.
- save\_only\_model (bool, optional, defaults to False) When checkpointing, whether to only save the model, or also the optimizer, scheduler & rng state. Note that when this is true, you won't be able to resume training from checkpoint. This enables you to save storage by not storing the optimizer, scheduler & rng state. You can only load the model using from\_pretrained with this option set to True.
- restore\_callback\_states\_from\_checkpoint (boo1, optional, defaults to False) Whether to restore the callback states from the checkpoint. If True, will override callbacks passed to the Trainer if they exist in the checkpoint."
- **use\_cpu** (bool, *optional*, defaults to False) Whether or not to use cpu. If set to False, we will use cuda or mps device if available.
- **seed** (int, *optional*, defaults to 42) Random seed that will be set at the beginning of training. To ensure reproducibility across runs, use the ~Trainer.model\_init function to instantiate the model if it has some randomly initialized parameters.
- data\_seed (int, optional) Random seed to be used with data samplers. If not set, random generators for data sampling will use the same seed as seed. This can be used to ensure reproducibility of data sampling, independent of the model seed.
- **jit\_mode\_eval** (bool, *optional*, defaults to False) Whether or not to use PyTorch jit trace for inference.
- **use\_ipex** (bool, *optional*, defaults to False) Use Intel extension for PyTorch when it is available. IPEX installation.
- **bf16** (bool, *optional*, defaults to False) Whether to use bf16 16-bit (mixed) precision training instead of 32-bit training. Requires Ampere or higher NVIDIA architecture or using CPU (use\_cpu) or Ascend NPU. This is an experimental API and it may change.
- **fp16** (bool, *optional*, defaults to False) Whether to use fp16 16-bit (mixed) precision training instead of 32-bit training.
- **fp16\_opt\_level** (str, optional, defaults to 'O1') For fp16 training, Apex AMP optimization level selected in ['O0', 'O1', 'O2', and 'O3']. See details on the <u>Apex documentation</u>.
- **fp16\_backend** (str, optional, defaults to "auto") This argument is deprecated. Use half\_precision\_backend instead.
- half\_precision\_backend (str, optional, defaults to "auto") The backend to use for mixed precision training. Must be one of "auto", "apex", "cpu\_amp". "auto" will use CPU/CUDA AMP or APEX depending on the PyTorch version detected, while the other choices will force the requested backend.

• **bf16\_full\_eval** (boo1, *optional*, defaults to False) — Whether to use full bfloat16 evaluation instead of 32-bit. This will be faster and save memory but can harm metric values. This is an experimental API and it may change.

- **fp16\_full\_eval** (boo1, *optional*, defaults to False) Whether to use full float16 evaluation instead of 32-bit. This will be faster and save memory but can harm metric values.
- tf32 (boo1, optional) Whether to enable the TF32 mode, available in Ampere and newer GPU architectures. The default value depends on PyTorch's version default of torch.backends.cuda.matmul.allow\_tf32. For more details please refer to the TF32 documentation. This is an experimental API and it may change.
- local\_rank (int, optional, defaults to -1) Rank of the process during distributed training.
- ddp\_backend (str, optional) The backend to use for distributed training. Must be one of "nccl", "mpi", "ccl", "gloo", "hccl".
- **tpu\_num\_cores** (int, *optional*) When training on TPU, the number of TPU cores (automatically passed by launcher script).
- **dataloader\_drop\_last** (bool, *optional*, defaults to False) Whether to drop the last incomplete batch (if the length of the dataset is not divisible by the batch size) or not.
- eval\_steps (int or float, optional) Number of update steps between two evaluations if eval\_strategy="steps". Will default to the same value as logging\_steps if not set. Should be an integer or a float in range [0,1). If smaller than 1, will be interpreted as ratio of total training steps.
- **dataloader\_num\_workers** (int, *optional*, defaults to 0) Number of subprocesses to use for data loading (PyTorch only). 0 means that the data will be loaded in the main process.
- past\_index (int, optional, defaults to -1) Some models like <u>TransformerXL</u> or <u>XLNet</u> can make use of the past hidden states for their predictions. If this argument is set to a positive int, the <u>Trainer</u> will use the corresponding output (usually index 2) as the past state and feed it to the model at the next training step under the keyword argument mems.
- run\_name (str, optional, defaults to output\_dir) A descriptor for the run. Typically used for wandb and mlflow logging. If not specified, will be the same as output\_dir.
- disable\_tqdm (bool, optional) Whether or not to disable the tqdm progress bars and table of metrics produced by ~notebook.NotebookTrainingTracker in Jupyter Notebooks. Will default to True if the logging level is set to warn or lower (default), False otherwise.
- **remove\_unused\_columns** (boo1, *optional*, defaults to True) Whether or not to automatically remove the columns unused by the model forward method.
- **label\_names** (List[str], *optional*) The list of keys in your dictionary of inputs that correspond to the labels.
  - Will eventually default to the list of argument names accepted by the model that contain the word "label", except if the model used is one of the XxxForQuestionAnswering in which case it will also include the ["start\_positions", "end\_positions"] keys.
- load\_best\_model\_at\_end (boo1, optional, defaults to False) Whether or not to load the best
  model found during training at the end of training. When this option is enabled, the best checkpoint

will always be saved. See <u>save total limit</u> for more.

When set to True, the parameters save\_strategy needs to be the same as eval\_strategy, and in the case it is "steps", save\_steps must be a round multiple of eval\_steps.

metric\_for\_best\_model (str, optional) — Use in conjunction with load\_best\_model\_at\_end to specify the metric to use to compare two different models. Must be the name of a metric returned by the evaluation with or without the prefix "eval\_". Will default to "loss" if unspecified and load\_best\_model\_at\_end=True (to use the evaluation loss).

If you set this value, greater\_is\_better will default to True. Don't forget to set it to False if your metric is better when lower.

- greater\_is\_better (bool, optional) Use in conjunction with load\_best\_model\_at\_end and metric\_for\_best\_model to specify if better models should have a greater metric or not. Will default to:
  - •True if metric\_for\_best\_model is set to a value that isn't "loss" or "eval\_loss".
  - False if metric\_for\_best\_model is not set, or set to "loss" or "eval\_loss".
- **ignore\_data\_skip** (bool, *optional*, defaults to False) When resuming training, whether or not to skip the epochs and batches to get the data loading at the same stage as in the previous training. If set to True, the training will begin faster (as that skipping step can take a long time) but will not yield the same results as the interrupted training would have.
- **fsdp** (bool, str or list of FSDPOption, *optional*, defaults to '') Use PyTorch Distributed Parallel Training (in distributed training only).

A list of options along the following:

- "full\_shard": Shard parameters, gradients and optimizer states.
- "shard\_grad\_op": Shard optimizer states and gradients.
- "hybrid\_shard": Apply FULL\_SHARD within a node, and replicate parameters across nodes.
- "hybrid\_shard\_zero2": Apply SHARD\_GRAD\_OP within a node, and replicate parameters across nodes.
- •"offload": Offload parameters and gradients to CPUs (only compatible with "full\_shard" and "shard\_grad\_op").
- "auto\_wrap": Automatically recursively wrap layers with FSDP using default\_auto\_wrap\_policy.
- **fsdp\_config** (str or dict, *optional*) Config to be used with fsdp (Pytorch Distributed Parallel Training). The value is either a location of fsdp json config file (e.g., fsdp\_config.json) or an already loaded json file as dict.

A List of config and its options:

min\_num\_params (int, optional, defaults to 0): FSDP's minimum number of parameters for Default Auto Wrapping. (useful only when fsdp field is passed).

- •transformer\_layer\_cls\_to\_wrap (List[str], optional): List of transformer layer class names (case-sensitive) to wrap, e.g, BertLayer, GPTJBlock, T5Block ... (useful only when fsdp flag is passed).
- backward\_prefetch (str, optional) FSDP's backward prefetch mode. Controls when to prefetch next set of parameters (useful only when fsdp field is passed).

A list of options along the following:

- "backward\_pre": Prefetches the next set of parameters before the current set of parameter's gradient computation.
- "backward\_post": This prefetches the next set of parameters after the current set of parameter's gradient computation.
- •forward\_prefetch (bool, *optional*, defaults to False) FSDP's forward prefetch mode (useful only when fsdp field is passed). If "True", then FSDP explicitly prefetches the next upcoming all-gather while executing in the forward pass.
- •limit\_all\_gathers (boo1, optional, defaults to False) FSDP's limit\_all\_gathers (useful only when fsdp field is passed). If "True", FSDP explicitly synchronizes the CPU thread to prevent too many in-flight all-gathers.
- •use\_orig\_params (bool, *optional*, defaults to True) If "True", allows non-uniform requires\_grad during init, which means support for interspersed frozen and trainable paramteres. Useful in cases such as parameter-efficient fine-tuning. Please refer this [blog] (<a href="https://dev-discuss.pytorch.org/t/rethinking-pytorch-fully-sharded-data-parallel-fsdp-from-first-principles/1019">https://dev-discuss.pytorch.org/t/rethinking-pytorch-fully-sharded-data-parallel-fsdp-from-first-principles/1019</a>
- •sync\_module\_states (bool, *optional*, defaults to True) If "True", each individually wrapped FSDP unit will broadcast module parameters from rank 0 to ensure they are the same across all ranks after initialization
- •cpu\_ram\_efficient\_loading (bool, optional, defaults to False) If "True", only the first process loads the pretrained model checkpoint while all other processes have empty weights. When this setting as "True", sync\_module\_states also must to be "True", otherwise all the processes except the main process would have random weights leading to unexpected behaviour during training.
- activation\_checkpointing (bool, optional, defaults to False): If "True", activation checkpointing is a technique to reduce memory usage by clearing activations of certain layers and recomputing them during a backward pass. Effectively, this trades extra computation time for reduced memory usage.
- •xla (bool, *optional*, defaults to False): Whether to use PyTorch/XLA Fully Sharded Data Parallel Training. This is an experimental feature and its API may evolve in the future.

xla\_fsdp\_settings (dict, optional) The value is a dictionary which stores the XLA FSDP wrapping parameters.

For a complete list of options, please see <u>here</u>.

- •xla\_fsdp\_grad\_ckpt (boo1, optional, defaults to False): Will use gradient checkpointing over each nested XLA FSDP wrapped layer. This setting can only be used when the xla flag is set to true, and an auto wrapping policy is specified through fsdp\_min\_num\_params or fsdp\_transformer\_layer\_cls\_to\_wrap.
- deepspeed (str or dict, optional) Use <u>Deepspeed</u>. This is an experimental feature and its API may evolve in the future. The value is either the location of DeepSpeed json config file (e.g., ds\_config.json) or an already loaded json file as a dict"

If enabling any Zero-init, make sure that your model is not initialized until \*after\* initializing the `TrainingArguments`, else it will not be applied.

accelerator\_config (str, dict, or AcceleratorConfig, optional) — Config to be used with the internal Accelerator implementation. The value is either a location of accelerator json config file (e.g., accelerator\_config.json), an already loaded json file as dict, or an instance of AcceleratorConfig.

A list of config and its options:

- •split\_batches (bool, optional, defaults to False): Whether or not the accelerator should split the batches yielded by the dataloaders across the devices. If True the actual batch size used will be the same on any kind of distributed processes, but it must be a round multiple of the num\_processes you are using. If False, actual batch size used will be the one set in your script multiplied by the number of processes.
- dispatch\_batches (bool, optional): If set to True, the dataloader prepared by the Accelerator is only iterated through on the main process and then the batches are split and broadcast to each process. Will default to True for DataLoader whose underlying dataset is an IterableDataset, False otherwise.
- even\_batches (bool, optional, defaults to True): If set to True, in cases where the total batch size across all processes does not exactly divide the dataset, samples at the start of the dataset will be duplicated so the batch can be divided equally among all workers.
- •use\_seedable\_sampler (bool, optional, defaults to True): Whether or not use a fully seedable random sampler (accelerate.data\_loader.SeedableRandomSampler). Ensures training results are fully reproducable using a different sampling technique. While seed-to-seed results may differ, on average the differences are neglible when using multiple different seeds to compare. Should also be ran with ~utils.set\_seed for the best results.
- label\_smoothing\_factor (float, optional, defaults to 0.0) The label smoothing factor to use. Zero means no label smoothing, otherwise the underlying onehot-encoded labels are changed from 0s and 1s to label\_smoothing\_factor/num\_labels and 1 label\_smoothing\_factor + label\_smoothing\_factor/num\_labels respectively.

**debug** (str or list of DebugOption, *optional*, defaults to "") — Enable one or more debug features. This is an experimental feature.

## Possible options are:

- "underflow\_overflow": detects overflow in model's input/outputs and reports the last frames that led to the event
- "tpu\_metrics\_debug": print debug metrics on TPU

The options should be separated by whitespaces.

- **optim** (str or training\_args.OptimizerNames, *optional*, defaults to "adamw\_torch") The optimizer to use: adamw\_hf, adamw\_torch, adamw\_torch\_fused, adamw\_apex\_fused, adamw\_anyprecision or adafactor.
- optim\_args (str, optional) Optional arguments that are supplied to AnyPrecisionAdamW.
- **group\_by\_length** (bool, *optional*, defaults to False) Whether or not to group together samples of roughly the same length in the training dataset (to minimize padding applied and be more efficient). Only useful if applying dynamic padding.
- length\_column\_name (str, optional, defaults to "length") Column name for precomputed lengths. If the column exists, grouping by length will use these values rather than computing them on train startup. Ignored unless group\_by\_length is True and the dataset is an instance of Dataset.
- report\_to (str or List[str], optional, defaults to "all") The list of integrations to report the results and logs to. Supported platforms are "azure\_ml", "clearml", "codecarbon", "comet\_ml", "dagshub", "dvclive", "flyte", "mlflow", "neptune", "tensorboard", and "wandb". Use "all" to report to all integrations installed, "none" for no integrations.
- **ddp\_find\_unused\_parameters** (bool, *optional*) When using distributed training, the value of the flag find\_unused\_parameters passed to DistributedDataParallel. Will default to False if gradient checkpointing is used, True otherwise.
- **ddp\_bucket\_cap\_mb** (int, *optional*) When using distributed training, the value of the flag bucket\_cap\_mb passed to DistributedDataParallel.
- **ddp\_broadcast\_buffers** (bool, *optional*) When using distributed training, the value of the flag broadcast\_buffers passed to DistributedDataParallel. Will default to False if gradient checkpointing is used, True otherwise.
- **dataloader\_pin\_memory** (bool, *optional*, defaults to True) Whether you want to pin memory in data loaders or not. Will default to True.
- dataloader\_persistent\_workers (boo1, optional, defaults to False) If True, the data loader will
  not shut down the worker processes after a dataset has been consumed once. This allows to
  maintain the workers Dataset instances alive. Can potentially speed up training, but will increase
  RAM usage. Will default to False.
- dataloader\_prefetch\_factor (int, optional) Number of batches loaded in advance by each worker.
   2 means there will be a total of 2 \* num\_workers batches prefetched across all workers.

• **skip\_memory\_metrics** (boo1, *optional*, defaults to True) — Whether to skip adding of memory profiler reports to metrics. This is skipped by default because it slows down the training and evaluation speed.

• **push\_to\_hub** (bool, *optional*, defaults to False) — Whether or not to push the model to the Hub every time the model is saved. If this is activated, output\_dir will begin a git directory synced with the repo (determined by hub\_model\_id) and the content will be pushed each time a save is triggered (depending on your save\_strategy). Calling <u>save\_model()</u> will also trigger a push.

If output\_dir exists, it needs to be a local clone of the repository to which the <u>Trainer</u> will be pushed.

- **resume\_from\_checkpoint** (str, optional) The path to a folder with a valid checkpoint for your model. This argument is not directly used by <u>Trainer</u>, it's intended to be used by your training/evaluation scripts instead. See the <u>example scripts</u> for more details.
- hub\_model\_id (str, optional) The name of the repository to keep in sync with the local output\_dir. It can be a simple model ID in which case the model will be pushed in your namespace. Otherwise it should be the whole repository name, for instance "user\_name/model", which allows you to push to an organization you are a member of with "organization\_name/model". Will default to user\_name/output\_dir\_name with output\_dir\_name being the name of output\_dir.

Will default to the name of output\_dir.

- hub\_strategy (str or HubStrategy, optional, defaults to "every\_save") Defines the scope of what is pushed to the Hub and when. Possible values are:
  - "end": push the model, its configuration, the tokenizer (if passed along to the <u>Trainer</u>) and a draft of a model card when the <u>save model()</u> method is called.
  - "every\_save": push the model, its configuration, the tokenizer (if passed along to the <u>Trainer</u>) and a draft of a model card each time there is a model save. The pushes are asynchronous to not block training, and in case the save are very frequent, a new push is only attempted if the previous one is finished. A last push is made with the final model at the end of training.
  - "checkpoint": like "every\_save" but the latest checkpoint is also pushed in a subfolder named last-checkpoint, allowing you to resume training easily with trainer.train(resume\_from\_checkpoint="last-checkpoint").
  - "all\_checkpoints": like "checkpoint" but all checkpoints are pushed like they appear in the output folder (so you will get one checkpoint folder per folder in your final repository)
- **hub\_token** (str, *optional*) The token to use to push the model to the Hub. Will default to the token in the cache folder obtained with huggingface-cli login.
- hub\_private\_repo (bool, optional, defaults to False) If True, the Hub repo will be set to private.
- hub\_always\_push (bool, optional, defaults to False) Unless this is True, the Trainer will skip pushing a checkpoint when the previous push is not finished.
- **gradient\_checkpointing** (bool, *optional*, defaults to False) If True, use gradient checkpointing to save memory at the expense of slower backward pass.

**gradient\_checkpointing\_kwargs** (dict, *optional*, defaults to None) — Key word arguments to be passed to the gradient\_checkpointing\_enable method.

- include\_inputs\_for\_metrics (bool, optional, defaults to False) Whether or not the inputs will be passed to the compute\_metrics function. This is intended for metrics that need inputs, predictions and references for scoring calculation in Metric class.
- eval\_do\_concat\_batches (boo1, optional, defaults to True) Whether to recursively concat inputs/losses/labels/predictions across batches. If False, will instead store them as lists, with each batch kept separate.
- auto\_find\_batch\_size (bool, optional, defaults to False) Whether to find a batch size that will fit
  into memory automatically through exponential decay, avoiding CUDA Out-of-Memory errors.
   Requires accelerate to be installed (pip install accelerate)
- **full\_determinism** (boo1, *optional*, defaults to False) If True, <u>enable full determinism()</u> is called instead of <u>set\_seed()</u> to ensure reproducible results in distributed training. Important: this will negatively impact the performance, so only use it for debugging.
- **torchdynamo** (str, *optional*) If set, the backend compiler for TorchDynamo. Possible choices are "eager", "aot\_eager", "inductor", "nvfuser", "aot\_nvfuser", "aot\_cudagraphs", "ofi", "fx2trt", "onnxrt" and "ipex".
- ray\_scope (str, optional, defaults to "last") The scope to use when doing hyperparameter search with Ray. By default, "last" will be used. Ray will then use the last checkpoint of all trials, compare those, and select the best one. However, other options are also available. See the <a href="Ray documentation">Ray documentation</a> for more options.
- ddp\_timeout (int, optional, defaults to 1800) The timeout for torch.distributed.init\_process\_group calls, used to avoid GPU socket timeouts when performing slow operations in distributed runnings. Please refer the [PyTorch documentation] (https://pytorch.org/docs/stable/distributed.html#torch.distributed.init\_process\_group) for more information.
- **use\_mps\_device** (bool, *optional*, defaults to False) This argument is deprecated.mps device will be used if it is available similar to cuda device.
- **torch\_compile** (bool, *optional*, defaults to False) Whether or not to compile the model using PyTorch 2.0 <u>torch.compile</u>.

This will use the best defaults for the <u>torch.compile API</u>. You can customize the defaults with the argument torch\_compile\_backend and torch\_compile\_mode but we don't guarantee any of them will work as the support is progressively rolled in in PyTorch.

This flag and the whole compile API is experimental and subject to change in future releases.

• **torch\_compile\_backend** (str, *optional*) — The backend to use in torch.compile. If set to any value, torch\_compile will be set to True.

Refer to the PyTorch doc for possible values and note that they may change across PyTorch versions.

This flag is experimental and subject to change in future releases.

• torch\_compile\_mode (str, optional) — The mode to use in torch.compile. If set to any value, torch\_compile will be set to True.

Refer to the PyTorch doc for possible values and note that they may change across PyTorch versions.

This flag is experimental and subject to change in future releases.

• **split\_batches** (bool, *optional*) — Whether or not the accelerator should split the batches yielded by the dataloaders across the devices during distributed training. If

set to True, the actual batch size used will be the same on any kind of distributed processes, but it must be a

round multiple of the number of processes you are using (such as GPUs).

• **include\_tokens\_per\_second** (bool, *optional*) — Whether or not to compute the number of tokens per second per device for training speed metrics.

This will iterate over the entire training dataloader once beforehand,

and will slow down the entire process.

• **include\_num\_input\_tokens\_seen** (bool, *optional*) — Whether or not to track the number of input tokens seen throughout training.

May be slower in distributed training as gather operations must be called.

- neftune\_noise\_alpha (Optional[float]) If not None, this will activate NEFTune noise embeddings. This can drastically improve model performance for instruction fine-tuning. Check out the <u>original paper</u> and the <u>original code</u>. Support transformers PreTrainedModel and also PeftModel from peft.
- optim\_target\_modules (Union[str, List[str]], optional) The target modules to optimize, i.e. the module names that you would like to train, right now this is used only for GaLore algorithm <a href="https://arxiv.org/abs/2403.03507">https://arxiv.org/abs/2403.03507</a> See: <a href="https://github.com/jiaweizzhao/GaLore">https://github.com/jiaweizzhao/GaLore</a> for more details. You need to make sure to pass a valid GaloRe optimizer, e.g. one of: "galore\_adamw", "galore\_adamw\_8bit", "galore\_adafactor" and make sure that the target modules are nn.Linear modules only.
- batch\_eval\_metrics (Optional[bool], defaults to False) If set to True, evaluation will call compute\_metrics at the end of each batch to accumulate statistics rather than saving all eval logits in memory. When set to True, you must pass a compute\_metrics function that takes a boolean argument compute\_result, which when passed True, will trigger the final global summary statistics from the batch-level summary statistics you've accumulated over the evaluation set.

TrainingArguments is the subset of the arguments we use in our example scripts **which relate to the training loop itself**.

Using <u>HfArgumentParser</u> we can turn this class into <u>argparse</u> arguments that can be specified on the command line.