```
import argparse
import collections
import logging
import math
import os
import re
import time
from concurrent.futures import ProcessPoolExecutor
import numpy as np
import smdistributed.modelparallel
import smdistributed.modelparallel.torch as smp
import torch
import torch.nn as nn
import torch.utils.data
import transformers
from data pipeline import create pretraining dataloader
from fp16 import FP16_Module, FP16_Optimizer, load_fp16_optimizer,
save fp16 optimizer
from learning_rates import AnnealingLR
from smdistributed.modelparallel.torch.nn import FusedLayerNorm as
LayerNorm
from smdistributed.modelparallel.torch.nn.huggingface.gpt2 import (
    translate hf state dict to smdistributed,
    translate state dict to hf gpt2,
from torch import optim
from torch.nn.parallel.distributed import DistributedDataParallel
from transformers import (
    CONFIG MAPPING,
    MODEL FOR CAUSAL LM MAPPING,
    AutoConfig,
    AutoModelForCausalLM,
    AutoTokenizer,
    GPT2Config,
    default data collator,
    set seed,
from transformers.trainer utils import is main process
logger = logging.getLogger( name )
def get learning rate scheduler (optimizer, args):
    # Add linear learning rate scheduler.
    if args.lr decay iters is not None:
        num_iters = args.lr_decay_iters
    else:
        num iters = args.max steps
    num iters = max(1, num iters)
    init step = 0
    warmup iter = args.warmup * num iters
    plateau iter = warmup iter + args.plateau * num iters
```

```
lr scheduler = AnnealingLR(
        optimizer,
        start lr=args.lr,
        warmup iter=warmup iter,
        plateau iter=plateau iter,
        total iters=num iters,
        decay style=args.lr decay style,
        last iter=init step,
        min lr=args.min lr,
        use checkpoint lr scheduler=args.load partial or args.load full,
        override lr scheduler=False,
    )
    return lr scheduler
def get param groups by weight decay(module):
    weight_decay_params = {"params": []}
    no weight decay params = {"params": [], "weight decay": 0.0}
    param ids = set()
    for module_ in module.modules():
        if isinstance(module , LayerNorm):
            for p in list(module . parameters.values()):
                if p is not None and id(p) not in param ids:
                    no_weight_decay_params["params"].append(p)
                    param ids.add(id(p))
        else:
            for n, p in list(module . parameters.items()):
                if p is not None and \overline{n} != "bias" and id(p) not in
param ids:
                    weight decay params["params"].append(p)
                    param ids.add(id(p))
            for n, p in list(module_._parameters.items()):
                if p is not None and n == "bias" and id(p) not in
param ids:
                    no weight decay params["params"].append(p)
                    param ids.add(id(p))
    return weight decay params, no weight decay params
# SMP modification: Define smp.step. Return any tensors needed outside.
@smp.step
def train step (model, optimizer, input ids, attention mask, args):
    if args.logits output:
        output = model(input ids=input ids,
attention mask=attention mask, labels=input ids)
        loss = output["loss"]
    else:
        loss = model(input ids=input ids, attention mask=attention mask,
labels=input ids)["loss"]
    if args.fp16:
        optimizer.backward(loss, update master grads=False)
    else:
        model.backward(loss)
```

```
if args.logits output:
        return output
    return loss
# SMP modification: Define smp.step. Return any tensors needed outside.
@smp.step
def test step (model, input ids, attention mask):
    loss = model(input ids=input ids, attention mask=attention mask,
labels=input ids)["loss"]
   return loss
def save ckptsum(args, model, optimizer, filename):
    results = collections.defaultdict(dict)
   model result = collections.defaultdict(dict)
    if args.fp16:
        from fp16.fp16util import register optimizer hooks
        register optimizer hooks (model)
    def get optimizer result(optimizer states):
        optimizer result = collections.defaultdict(dict)
        for param idx, state in optimizer states.items():
            for key, val in state.items():
                if isinstance(val, torch.Tensor):
                    optimizer result["tensors"][f"{param idx} {key}"] =
torch.sum(val)
                else:
                    optimizer result["scalars"][f"{param idx} {key}"] =
val
        return optimizer result
    if not args.shard optimizer state:
        optimizer result =
get optimizer result(optimizer.local state dict()["state"])
        local state dict = optimizer.local state dict()["state"]
        if smp.rdp rank() == 0:
            optimizer result = []
            for partial local state dict in local state dict:
optimizer result.append( get optimizer result(partial local state dict))
    for param name, param in model.local state dict().items():
        if isinstance(param, torch.Tensor):
            model result["tensors"][param name] = torch.sum(param)
        else:
            model result["scalars"][param name] = param
    if smp.rdp rank() == 0:
        results["optimizer"] = optimizer result
        results["model"] = model result
```

```
def load and verify ckptsum(args, model, optimizer, filename):
    results = smp.load(filename)
    optimizer result = (
        results["optimizer"]
        if not args.shard optimizer state
        else results["optimizer"][smp.rdp rank()]
    model result = results["model"]
    def opt check fn(mod, opt):
        loaded opt states = (
            opt.orig state dict()["state"]
            if args.shard optimizer state
            else opt.local state dict()["state"]
        for param idx, state in loaded opt states.items():
            for key, val in state.items():
                if isinstance(val, torch.Tensor):
                    assert torch.isclose(
                        torch.sum(val),
optimizer_result["tensors"][f"{param_idx}_{key}"]
                    ), f"mismatch for param idx: {param idx}, key is
{key}"
                else:
                    assert (
                        val ==
optimizer result["scalars"][f"{param idx} {key}"]
                    ), f"mismatch for param idx: {param idx}, key is
{ key} "
        print("Optimizer save/load check passed successfully")
    def model check fn(mod, opt):
        for param name, param in mod.local state dict().items():
            if isinstance(param, torch.Tensor):
                assert torch.isclose(
                    torch.sum(param), model result["tensors"][param_name]
                ), f"mismatch for param name: {param name}"
            else:
                assert (
                    param == model result["scalars"][param name]
                ), f"mismatch for param name: {param name}"
        print("Model save/load check passed successfully")
    model.register post partition hook (model check fn)
    model.register post step hook(opt check fn)
def save(
   output save file,
   model,
    optimizer,
```

smp.save(results, filename)

```
lr scheduler,
    model config,
    num params,
    total steps,
    curr train path index,
    args,
    partial=True,
    translate to hf=False,
    seq length=1024,
    batch idx=0,
):
    save_fn = save_fp16_optimizer
    save dict = {
        "cli args": args.__dict__,
        "num params": num params,
        "total steps": total steps,
        "curr train path index": curr train path index,
        "model_config": model_config,
        "batch idx": batch idx,
    }
    if lr scheduler is not None:
        save dict["lr scheduler"] = lr scheduler.state dict()
    if partial:
        # SMP modification: check if using optimizer state sharding or
tensor parallelism
        if args.gather if shard > 0 or smp.rdp rank() == 0:
            # if not gather the opt checkpoint, only save the model for
rdp rank 0
            save dict["model"] = model.local state dict()
    else:
        model state dict = model.state dict(gather to rank0=True)
        if smp.rank() == 0:
            save dict["model"] = (
                translate state dict to hf gpt2 (model state dict,
seq length)
                if translate to hf
                else model state dict
            )
    if args.fp16:
        if not partial and args.skip full optimizer:
            print("Skipping saving the final optimizer state")
            if args.shard optimizer state == 0 or partial:
                save dict["optimizer"] = save fn(args, model, optimizer,
partial=partial)
            else:
                     "Saving the full optimizer state does not work with
shard optimizer state > 0! Skipping..."
    else:
        # fp32
```

```
if partial:
            save dict["optimizer"] = optimizer.local state dict()
            if not args.skip full optimizer:
                save dict["optimizer"] = optimizer.state dict()
            else:
                print("Skipping saving of full optimizer state")
    # SMP modification: criteria for checkpointing the zeroth rank for
    # pipeline parallelism, checkpointing the zeroth reduced data
parallel
   # rank for tensor parallelism, and preventing checkpointing if
optimizer
    # state sharding is enabled
    if not args.gather if shard or (smp.rdp rank() == 0 and partial) or
smp.rank() == 0:
        smp.save(save dict, output save file, partial=partial, v3=not
args.gather if shard)
    print(f"Finished checkpointing after {total steps} steps:
{output save file}")
def load model and optimizer (
    output dir,
    model,
    optimizer,
    lr scheduler,
    partial,
    args,
    translate from hf=False,
    seq_length=102\overline{4},
    load model=True,
    load optimizer=True,
    num params=0,
):
    # Find longest-trained checkpoint
    re pattern = f"trained_gpt_nparams-{num_params}_steps-
(?P<total steps>\d+)\.pt"
    if partial:
        re_pattern += "_(?P<rank>\d+)"
    else:
        re pattern += "$"
    ckpt paths = sorted(
            (int(re.match(re pattern, p).group("total steps")),
os.path.join(output dir, p))
            for p in os.listdir(output dir)
            if re.match(re pattern, p)
        reverse=True,
    if not ckpt paths:
```

```
raise Exception (
            f'No checkpoints could be found in "{output dir}".
Candidates: {os.listdir(output dir)}'
    local ckpt path = ckpt paths[0][1]
    if partial:
        # need to pass prefix without ranks to smp
        local ckpt path = local ckpt path.split(".pt")[0] + ".pt"
    if args.gather if shard > 0:
        # Should expect v2 checkpoint here
        checkpoint = smp.load(local ckpt path, partial=partial)
    else:
        # Loading separately for model and opt
        checkpoint =
torch.load(f"{local_ckpt_path}_{smp.pp_rank()}_{smp.tp_rank()}_0")
        if smp.rdp rank() != 0:
            opt checkpoint = torch.load(
f"{local ckpt path} {smp.pp rank()} {smp.tp rank()} {smp.rdp rank()}"
    if load model:
        checkpointed model = (
            translate hf state dict to smdistributed(checkpoint["model"],
seq length)
            if translate from hf
            else checkpoint["model"]
        model.load state dict(checkpointed model,
same partition load=args.same partition load > 0)
        if lr scheduler is not None:
            lr scheduler.load state dict(checkpoint["lr scheduler"])
    if load optimizer:
        # Loading loss scale eagerly
        opt state dict = checkpoint["optimizer"]
        optimizer.loss scaler = opt state dict["loss scaler"]
        optimizer.loss scaler.model = model
        optimizer.dynamic loss scale =
opt state dict["dynamic loss scale"]
        optimizer.overflow = opt state dict["overflow"]
        optimizer.first closure call this step =
opt state dict["first closure call this step"]
        def opt load hook(mod, opt):
            load fn = load fp16 optimizer
            if args.fp16:
                if not partial and args.skip full optimizer:
                    print(
                        "Skipping loading the final optimizer state, and
reloading master params from model params"
```

```
opt.reload model params()
                else:
                    load fn(args, mod, opt, checkpoint, partial=partial)
            else:
                # fp32
                if not partial and args.skip full optimizer:
                    print ("Skipping loading the final optimizer state")
                else:
                    opt.load state dict(checkpoint["optimizer"])
        model.register_post_step_hook(opt load hook)
    print(f'Loaded model from "{local ckpt path}"')
   batch idx = 0
    if "batch idx" in checkpoint:
        batch idx = checkpoint["batch idx"]
    return (
        model,
        optimizer,
        checkpoint["total steps"],
        checkpoint["curr_train_path_index"],
       batch idx,
    )
def delete oldest ckpt(args, delete on rank0 only=False):
    to delete = smp.rank() == 0 if delete on rank0 only else
smp.local rank() == 0
    if to delete:
        re pattern = "trained qpt nparams-(?P<num params>\d+) steps-
(?P<total steps>\d+)\.pt"
        # partial
        re pattern += " (?P < pp rank > d+) (?P < tp rank > d+)"
        paths per step = collections.defaultdict(list)
        for p in os.listdir(args.checkpoint dir):
            if re.match(re pattern, p):
                step = int(re.match(re pattern, p).group("total steps"))
                path = os.path.join(args.checkpoint dir, p)
                paths per step[step].append(path)
        if paths per step:
            oldest step = sorted(paths per step.keys())[0]
            num parts = len(paths per step[oldest step])
        if len(paths per step) > args.num kept checkpoints:
            # delete oldest step
                for p in paths per step[oldest step]:
                    os.remove(p)
```

```
# else We still haven't reached maximum number of checkpoints --
no need to delete older ones
    return None
def eval model (model, dataloader, num batches, use wiki data):
    model = model.eval()
    n \text{ batches} = 0
    loss = 0.0
    with torch.no grad():
        for batch_idx, input_data in enumerate(dataloader):
            if use wiki data:
                input ids, , attention mask, , = input data
            else:
                input ids, attention mask = input data
            if batch idx >= num_batches:
                break
            loss += test step(model, input ids,
attention mask).reduce mean()
            n batches += 1
    if n batches > 0:
        torch.distributed.all reduce(loss,
group=smp.get_dp_process_group())
        loss /= smp.dp size()
        loss /= n batches
        loss = loss.item()
        ppl = math.exp(loss)
    else:
        loss = -1.0
        ppl = -1.0
    return loss, ppl
def train(
   model,
    optimizer,
    lr scheduler,
    model_config,
    start train path index,
    start batch index,
    num params,
    total steps,
    args,
):
    model.train()
    if args.parallel proc data processing:
        pool = ProcessPoolExecutor(1)
    dp rank = smp.dp rank() if not args.prescaled batch else
smp.rdp_rank()
```

```
dp size = smp.dp size() if not args.prescaled batch else
smp.rdp size()
    data type = "wiki" if args.use wiki data else "openwebtext"
    if args.use wiki data:
        train paths = sorted(
                os.path.join(args.training dir, p)
                for p in os.listdir(args.training dir)
                if os.path.isfile(os.path.join(args.training dir, p)) and
"training" in p
        )
    else:
        if args.zipped data > 0:
            file extension = ".json.gz"
        else:
            file extension = ".json"
        train paths = sorted(
            Γ
                os.path.join(args.training dir, p)
                for p in os.listdir(args.training dir)
                if p.endswith(file extension)
            ]
        )
    train dataloader = create pretraining dataloader (
        [train paths[start train path index]],
        args.train batch size,
        args.max context width,
        seed=args.seed,
        dp rank=dp rank,
        dp size=dp size,
        shuffle=args.same seed < 1,</pre>
        zipped=args.zipped data > 0,
        use last file only=args.fast validation > 0,
        data type=data type,
    )
    if args.validation freq is not None:
        # load all validation examples
        if smp.rank() == 0:
            print("Creating val dataloader")
        if args.use wiki data:
            val paths = sorted(
                    os.path.join(args.test dir, p)
                    for p in os.listdir(args.test dir)
                    if os.path.isfile(os.path.join(args.test dir, p)) and
"testing" in p
                1
            )
        else:
```

```
if args.zipped data > 0:
                file extension = ".json.gz"
            else:
                file extension = ".json"
            val_paths = sorted(
                    os.path.join(args.test dir, p)
                    for p in os.listdir(args.test dir)
                    if p.endswith(file extension)
            )
        val_dataloader = create_pretraining_dataloader(
            val paths,
            args.val batch size,
            args.max context width,
            seed=args.seed,
            dp rank=dp rank,
            dp size=dp_size,
            shuffle=True,
            zipped=args.zipped data > 0,
            use last file only=args.fast validation > 0,
            data type=data type,
        )
        if smp.rank() == 0:
            print("Created val dataloader")
    start = time.time()
    throughput = None
    to_save = {"loss": [], "val_loss": []}
    loss\ metric = 0
    def should record():
        # only record the ranks that in the tp group that contains global
rank 0
        if smp.tp size() > 1:
            tp group = smp.get tp group()
            return 0 in tp group
        else:
            return smp.rank() == 0
    # Set the same seed for computation
    set seed(args.seed)
    for index in range(start train path index, args.epochs *
len(train paths)):
        next train path index = (index + 1) % len(train paths)
        curr train path index = index % len(train paths)
        if total steps >= args.max steps:
            break
        if args.parallel proc data processing:
            dataset future = pool.submit(
                create_pretraining_dataloader,
```

```
[train paths[next train path index]],
                args.train batch size,
                args.max context width,
                seed=args.seed,
                dp rank=dp rank,
                dp size=dp size,
                shuffle=args.same seed < 1,</pre>
                zipped=args.zipped data > 0,
                use last file only=args.fast validation > 0,
                data type=data type,
            )
        if smp.rank() == 0:
            if args.use wiki data:
                print(f"Reading data from training path
{train dataloader.dataset.input file}")
            else:
                print(f"Reading data from training path
{train_dataloader.dataset.input_paths}")
        for batch idx, input data in enumerate(train dataloader):
            if batch idx < start batch index:</pre>
                if smp.rank() == 0:
                    print(
                        f"Resuming from saved batch index
{start batch index}, skipping batch {batch idx}..."
                if start batch index == len(train dataloader):
                    # If saving at the last batch of the file, read from
the next file
                    start batch index = 0
                    break
                continue
            else:
                start batch index = 0
            if args.use wiki data:
                input ids, , attention mask, , = input data
            else:
                input ids, attention mask = input data
            if total_steps >= args.max_steps:
                break
            step start = time.time()
            if args.fp16:
                optimizer.zero_grad(set_grads_to_None=True)
            else:
                optimizer.zero grad()
            if args.logits output:
                train output = train step(model, optimizer, input ids,
attention mask, args)
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loss mb = train output["loss"]
                logits mb = train output["logits"]
                if smp.tp size() > 1:
                    logits = torch.cat(tuple(logits mb.outputs), dim=1)
                else:
                    logits = torch.cat(tuple(logits mb.outputs), dim=0)
            else:
                # Return value, loss mb is a StepOutput object
                loss mb = train step(model, optimizer, input ids,
attention mask, args)
            # SMP modification: Average the loss across microbatches.
            loss = loss mb.reduce mean()
            if not args.validation freq:
                loss metric = loss.item()
            if args.fp16:
                optimizer.update master grads()
                optimizer.clip master grads(args.grad clip)
                optimizer.step()
                overflow = optimizer.overflow
            else:
                optimizer.step()
            if not (args.fp16 and overflow):
                lr scheduler.step()
            if args.enable memory profiling > 0:
                memory status(msg="After opt step")
            total steps += 1
            time elapsed = time.time() - start
            step time = time.time() - step start
            sample processed = input ids.shape[0] * dp size
            throughput = sample processed / step time
            if smp.rank() == 0 and not total steps % args.logging freq:
                print(
                    f"({int(time elapsed)}s), Batch {total steps - 1}
Loss: {loss.item()}, Speed: {throughput} samples/sec"
            # evaluate on validation
            if args.validation freq and not (total steps %
args.validation freq):
                cur state = np.random.get state()
                model = model.eval()
                val loss, val ppl = eval model(
                    model, val dataloader, args.validation batches,
args.use wiki data
                if is main process(smp.rank()):
                    print(
                        f"({int(time.time()-start)}s) Batch {total steps
- 1} Validation loss: {val_loss}"
```

```
)
                    print(
                         f"({int(time.time()-start)}s) Batch {total steps
- 1} Validation perplexity: {val ppl}"
                loss metric = val loss
                if args.logits output:
                    to save["val loss"].append(val loss)
                model = model.train()
                if args.preserve np state > 0:
                    np.random.set state(cur state)
            # checkpoint
            if not (total steps % args.checkpoint freq):
                base path = f"trained gpt nparams-{num params} steps-
{total steps}.pt"
                out path = os.path.join(args.checkpoint dir, base path)
                total_ckpts = total_steps // args.checkpoint_freq
                delete oldest ckpt(args,
delete on rank0 only=args.use fsx > 0)
                # save or verify ckptsum if this is the last checkpoint
                if (args.save_or_verify_ckptsum and total_steps >=
args.max steps) or (
                    (total_ckpts + 1) * args.checkpoint freq
                ) > args.max steps:
                    # Save optimizer and model tensor sums and scalars
before saving
                    save ckptsum(
                        args,
                        model,
                        optimizer,
                        filename=os.path.join(args.model dir,
"saved partial_sum"),
                save(
                    out path,
                    model,
                    optimizer,
                    lr scheduler,
                    model config,
                    num params,
                    total steps,
                    curr train_path_index,
                    args,
                    partial=True,
                    batch idx=batch idx + 1,
                )
            if args.logits output:
                to save["loss"].append(loss.item())
```

```
if total steps >= args.max steps:
            if should record() and args.logits output:
                to save["logits"] = logits.detach().cpu()
                output file = f"rank {smp.rank()} " + args.logits output
                torch.save(to save, os.path.join(args.model dir,
output file))
                print(f"logits and loss saved at
{os.path.join(args.model dir, output file)}")
            break
        del train dataloader
        if args.parallel proc data processing:
            s = time.time()
            train dataloader = dataset future.result(timeout=None)
            wait time = time.time() - s
            if wait time > 1:
                # TODO if this happens, we should try num workers>1 in
dataloader
                print(
                    f"[{smp.rank()}] Waited {wait time} for data loader
to be ready. Please check if dataloader performance can be improved to
avoid these waits."
        else:
            train dataloader = create pretraining dataloader(
                [train paths[next train path index]],
                args.train batch size,
                args.max context width,
                seed=args.seed,
                dp rank=dp rank,
                dp size=dp size,
                shuffle=args.same seed < 1,
                zipped=args.zipped data > 0,
                use last file only=args.fast validation > 0,
                data type=data type,
            )
    return total steps, throughput, loss metric
def parse args():
   parser = argparse.ArgumentParser()
    # hyperparameters sent by the client are passed as command-line
arguments to the script.
    opt grp = parser.add argument group(
        title="optimization", description="arguments for optimization"
    opt grp.add argument(
       "--train batch size",
        type=int,
        default=4,
```

```
help="batch size per dp rank, for tensor parallelism degree 8
with pipeline parallel degree 1 this means 8*this batch size per node",
    opt grp.add argument("--val batch size", type=int, default=4)
    opt grp.add argument("--max steps", type=int, default=5000)
    opt grp.add argument("--seed", type=int, default=12345)
    opt grp.add argument("--same seed", type=int, default=0)
    opt grp.add argument("--n gpus", type=str,
default=os.environ["SM NUM GPUS"])
    opt grp.add argument("--fp16", default=0, type=int, help="automatic
mixed precision training")
    opt grp.add argument(
        "--fp32 grad accumulation", default=0, type=int, help="Enable
FP32 Grad accumulation"
    opt grp.add argument("--grad clip", default=1.0, type=float,
help="gradient clipping")
    opt grp.add argument("--weight decay", default=0.01, type=float,
help="weight decay")
    opt grp.add argument(
        "--beta1", default=0.9, type=float, help="beta1 parameter for
Adam optimizer"
    )
    opt_grp.add_argument(
        "--beta\overline{2}", default=0.95, type=float, help="beta2 parameter for
Adam optimizer"
   )
    opt grp.add argument(
        "--activation checkpointing",
        type=int,
        default=1,
        help="enable gradient checkpointing to reduce memory
consumption",
    )
    parser.add argument (
        "--logging freq", type=int, default=1, help="number of iterations
between logging"
    # I/O
    io grp = parser.add argument group(title="io", description="location
for input and output")
    io grp.add argument ("--use wiki data", type=int, default=0, help="use
wiki corpus data for training")
    io grp.add argument ("--zipped data", type=int, default=1, help="input
data is zipped files")
    io grp.add argument(
        "--epochs", type=int, default=3, help="times of iterating over
the training dataset"
    io grp.add argument ("--output-data-dir", type=str,
default=os.environ["SM OUTPUT DATA DIR"])
    io grp.add argument (
        "--checkpoint-dir",
```

```
type=str,
        default="/opt/ml/checkpoints",
        help="Saves partial checkpoints (model, optimizer) to this dir,
and loads latest checkpoint from this if load partial is specified.",
    io grp.add argument(
        "--model-dir",
        type=str,
        default=os.environ["SM MODEL DIR"],
        help="Saves full model for inference to this dir. Also used if
load full is given to load the model. Note the lack of optimizer state
here.",
    io grp.add argument ("--training-dir", type=str,
default=os.environ["SM CHANNEL TRAIN"])
    io grp.add argument ("--test-dir", type=str,
default=os.environ["SM CHANNEL TEST"])
    io_grp.add argument(
        "--parallel proc data_processing",
        type=int,
        default=0,
        help="Load data in parallel with a different process. At any
point a process can have two files in memory. With tensor parallelism,
each of the 8 processes on an instance will then have 2 files in memory.
Depending on file sizes this may or may not be feasible. With pipeline
parallelism this was not a problem as only 1 rank on an instance loaded
data.",
    io grp.add argument(
        "--save final full model",
        type=int,
        default=0,
        help="Enabling this will save a combined model only at the end",
    io_grp.add argument(
        "--skip full optimizer",
        type=int,
        default=1,
        help="Disabling this will also save the full optimizer state",
    io grp.add argument ("--load partial", type=int, default=0, help="Load
from partial checkpoints")
    io grp.add argument("--load full", type=int, default=0, help="Load
from full checkpoints")
    io grp.add argument(
        "--logits output", type=str, default="", help="Path to save
    io grp.add argument("--prescaled batch", type=int, default=1,
help="use prescaled batch")
    # configure model size
    model grp = parser.add argument group(
```

```
title="model", description="arguments to describe model
configuration"
    model grp.add argument("--max context width", type=int, default=1024)
    model grp.add argument("--vocab size", type=int, default=50257)
    model grp.add argument("--hidden width", type=int, default=768)
    model grp.add argument("--num layers", type=int, default=12)
    model grp.add argument("--num heads", type=int, default=12)
    model grp.add argument("--resid pdrop", type=float, default=0.1)
    model_grp.add_argument("--embd_pdrop", type=float, default=0.1)
    model grp.add argument("--attn pdrop", type=float, default=0.1)
    model grp.add argument("--summary first pdrop", type=float,
default=0.1)
    model grp.add argument("--use adamw", type=int, default=0, help="Use
adamw optimizer")
    smp grp = parser.add argument group(title="smp", description="smp")
    smp_grp.add_argument("--tensor_parallel_degree", type=int, default=8)
    smp grp.add argument("--pipeline parallel degree", type=int,
default=1)
    smp grp.add argument("--microbatches", type=int, default=1)
    smp grp.add argument("--active microbatches", type=int, default=None)
    smp grp.add argument("--optimize", type=str, default="speed")
    smp grp.add argument("--activation strategy", type=str,
default="each")
    smp grp.add argument("--shard optimizer state", type=int, default=0)
    smp grp.add argument("--offload activations", type=int, default=0)
    smp grp.add argument("--fast mode", type=int, default=0)
    smp_grp.add_argument("--static mode", type=int, default=0)
    smp grp.add argument("--delayed param", type=int, default=0)
    smp grp.add argument("--same partition load", type=int, default=0)
    smp grp.add argument("--attention in fp32", type=int, default=0)
    smp_grp.add_argument("--placement strategy", type=str,
default="cluster")
    smp grp.add argument("--activation loading horizon", type=int,
default=4)
    smp grp.add argument("--skip tracing", type=int, default=0)
    smp grp.add argument("--query key layer scaling", type=int,
default=1)
    smp_grp.add_argument("--fused softmax", type=int, default=1)
    smp grp.add argument("--fused bias gelu", type=int, default=1)
    parser.add argument(
        "--num kept checkpoints",
        type=int,
        default=5,
        help="how many checkpoints to keep before deleting",
    parser.add argument(
        "--checkpoint freq",
        type=int,
        default=10000,
        help="number of iterations between checkpointing",
    )
```

```
parser.add argument(
        "--validation freq",
        type=int,
        default=None,
        help="number of iterations to print validation loss",
    )
    parser.add argument(
        "--validation batches",
        type=int,
        default=10,
        help="number of batches to estimate validation loss",
    parser.add argument(
        "--manual partition",
        type=int,
        default=0,
        help="evenly distribute layers across the partitions",
    parser.add argument(
        "--match weights", type=int, default=0, help="Get weights from
the original model"
    parser.add argument(
        "--preserve np state",
        type=int,
        default=0,
        help="Perserve the numpy random state between validation",
    parser.add argument(
        "--fast validation",
        type=int,
        default=1,
        help="Running validation only with the last data file for faster
speed",
    parser.add argument(
        "--gather if shard",
        type=int,
        default=1,
        help="When sharding opt states is enabled, gather the opt
checkpoint to rdp rank 0 during saving",
    parser.add argument(
        "--clean cache",
        type=int,
        default=0,
        help="Clean torch reserved memory at he end of every step",
    parser.add argument("--use fsx", type=int, default=0, help="Using FSx
for checkpointing")
    parser.add argument(
        "--enable memory profiling", type=int, default=0, help="Enable
memory profile"
    )
```

```
# learning rate
    lr grp = parser.add argument group(
        title="lr", description="arguments for learning rate schedule"
    lr grp.add argument("--lr", type=float, default=None, help="Initial
learning rate.")
    lr grp.add argument(
        "--lr decay style",
        type=str,
        default="linear",
        choices=["constant", "linear", "cosine", "exponential",
"plateau"],
        help="Learning rate decay function.",
    lr_grp.add argument(
        "--lr decay iters",
        type=int,
        default=None,
        help="number of iterations to decay learning rate over," " If
None defaults to train iters",
    lr grp.add argument(
        "--min lr",
        type=float,
        default=0.0,
        help="Minumum value for learning rate. The scheduler" "clip
values below this threshold.",
    lr grp.add argument(
        "--warmup",
        type=float,
        default=0.01,
        help="Percentage of total iterations to warmup on "
        "(.01 = 1 percent of all training iters).",
    lr grp.add argument(
        "--plateau",
        type=float,
        default=0.4,
        help="Percentage of total iterations to keep at max if using
plateau lr",
    )
    ci grp = parser.add argument group(title="ci", description="ci
related settings")
    ci grp.add argument ("--ci", default=False, action="store true",
help="Whether enable ci")
    ci grp.add argument("--time to train", type=int, help="time to train
threshold")
    ci grp.add argument("--throughput", type=float, help="throughput
threshold")
    ci grp.add argument("--loss", type=float, help="loss threshold")
    ci grp.add argument(
```

```
"--save or verify ckptsum", default=False, action="store true",
help="Whether to save sum"
    args, _ = parser.parse_known args()
    return args
def main():
    args = parse args()
    if args.shard optimizer state > 0 and not args.skip full optimizer:
        raise ValueError(
            "If shard optimizer state is enabled, skip full optimizer
must also be enabled. Full optimizer saving is currently not supported
under optimizer state sharding."
    # any value here is overriden by the config set in notebook when
launching the sagemaker job
    smp_config = {
        "ddp": True,
        "tensor parallel degree": args.tensor parallel degree,
        "pipeline parallel degree": args.pipeline parallel degree,
        "microbatches": args.microbatches,
        # if activation checkpointing true checkpoints transformer layers
below
        "checkpoint attentions": False if args.activation checkpointing
else True,
        "shard optimizer state": args.shard optimizer state > 0,
        "prescaled batch": args.prescaled batch > 0,
        " match weights": args.match weights > 0,
        "fp16 params": args.fp16 > 0,
        "offload activations": args.offload activations > 0,
        "optimize": args.optimize,
        "placement strategy": args.placement strategy,
        "activation loading horizon": args.activation_loading_horizon,
        "skip tracing": args.skip tracing > 0,
        "auto partition": False if args.manual partition else True,
        "default partition": 0,
        " fp32 grad accumulation": args.fp32 grad accumulation > 0,
        "static mode": args.static mode > 0,
        "fast mode": args.fast mode > 0,
    if args.active microbatches is not None:
        smp config["active microbatches"] = args.active microbatches
    smp.init(smp config) # initialize modelparallel library
    if smp.rank() == 0:
        print("Arguments:", args. dict )
        print(f"Transformers version: {transformers. version }")
        print(f"smdistributed.modelparallel version:
{smdistributed.modelparallel. version }")
```

```
print(f"smdistributed config: {smp config}")
    if args.save final full model and smp.rank() == 0:
        print(
            f"[Warning] Note that save final full model only saves the
final model at the end of all steps. It does not save optimizer state.
Optimizer state is only saved with partial models which are saved at
checkpointing freq during training. If you want to restart training you
need partial checkpoints."
    if smp.local rank() == 0:
        for path in [args.model dir, args.checkpoint dir]:
            if not os.path.exists(path):
                os.makedirs(path, exist ok=True)
    model config = GPT2Config(
        vocab size=args.vocab size,
        n positions=args.max context width,
        n embd=args.hidden width,
        n layer=args.num layers,
        n head=args.num heads,
        n inner=None,
        activation function="gelu new",
        resid pdrop=args.resid pdrop,
        embd pdrop=args.embd pdrop,
        attn pdrop=args.attn pdrop,
        layer norm epsilon=1e-05,
        initializer range=0.02,
        summary type="cls index",
        summary use proj=True,
        summary activation=None,
        summary_proj to labels=True,
        summary first dropout=args.summary first pdrop,
        # gradient checkpointing=args.gradient checkpointing > 0,
        use cache=False,
        bos token id=50256,
        eos token id=50256,
        return dict=True,
    )
    # the following improves start-up time by skipping proper
initialization
    # of weights in the original model. this is not a problem because
DistributedModel
    # will override those weights anyway when tensor parallel degree > 1.
    if smp.tp size() > 1 and args.match weights < 1:
        from transformers.modeling utils import PreTrainedModel
        PreTrainedModel.init weights = lambda x: None
    set seed(args.seed)
    if args.fp16:
```

```
torch.set default dtype(torch.float16)
    with smp.tensor parallelism(
        enabled=smp.tp size() > 1,
        attention in fp32=args.attention in fp32 > 0,
        query key layer_scaling=args.query_key_layer_scaling > 0,
        fused softmax=args.fused softmax > 0,
        fused bias gelu=args.fused bias gelu > 0,
    ):
        with smp.delay param initialization (
            enabled=(smp.tp size() > 1 and args.match weights < 1 and</pre>
args.delayed param > 0)
        ):
            model = AutoModelForCausalLM.from config(model config)
    torch.set default dtype(torch.float32)
    if args.fp16:
        model = FP16_Module(model)
    num_params = sum([np.prod(p.size()) for p in model.parameters()])
    if smp.rank() == 0:
        print(f"# total parameters: {num params}")
    # SMP modification: Set the device to the GPU ID used by the current
process.
    # Input tensors should be transferred to this device.
    torch.cuda.set_device(smp.local rank()) # restrict each process to its own device
    device = torch.device("cuda")
    if not args.same seed:
        # Set seed by tp rank to prevent weights from being the same on
different tp ranks
        set seed(args.seed + smp.tp rank())
    # SMP modification: Use the DistributedModel container to provide the
model
    # to be partitioned across different ranks. For the rest of the
script,
    # the returned DistributedModel object should be used in place of
    # the model provided for DistributedModel class instantiation.
    if args.fp16:
        torch.set default dtype(torch.float16)
    model = smp.DistributedModel (model, trace device="gpu") # wrap the model with smp.DistributedModel
    if args.fp16:
        m = model.module
    else:
        m = model
    if smp.tp size() > 1:
        transformer layers = m.module.module.transformer.seg layers
    else:
        transformer layers = m.module.module.transformer.h
```

```
if args.manual partition:
        print(f"Manual partition enabled")
        # evenly distribute layers across all partitions
        div, rem = divmod(args.num layers, smp.pp size())
        get num layers = lambda x: (div + 1 if x >= smp.pp size() - rem
else div)
        assignments = []
        for pp rank in range(smp.pp size()):
            nl = get num layers(pp rank)
            print(f"{nl} layers assigned to partition {pp rank}")
            assignments += [pp rank for in range(nl)]
        for i, c in enumerate(transformer layers.children()):
            smp.set partition(c, assignments[i])
    torch.set default dtype(torch.float32)
    iter model = model
    # Build parameter groups (weight decay and non-decay).
    while isinstance(iter model, (DistributedDataParallel, FP16 Module)):
        iter model = iter model.module
    param groups = get param groups by weight decay(iter model)
    if args.use adamw > 0:
        optimizer = optim.AdamW(
            param groups, betas=(args.beta1, args.beta2), lr=args.lr,
weight decay=args.weight decay
    else:
        optimizer = optim.Adam(
            param groups, betas=(args.beta1, args.beta2), lr=args.lr,
weight decay=args.weight decay
    if args.activation checkpointing:
        kwargs = {}
        if isinstance(transformer layers, nn.Sequential):
            kwargs["pack args as tuple"] = True
            kwargs["strategy"] = args.activation strategy
            smp.set activation checkpointing(transformer layers,
**kwargs)
        else:
            for c in transformer layers.children():
                smp.set activation checkpointing(c)
    if args.fp16:
        optimizer = FP16 Optimizer(
            model,
            optimizer,
            static loss scale=None,
            dynamic loss scale=True,
            use smp=True,
            dynamic loss args={"scale window": 1000, "min scale": 1,
"delayed shift": 2},
```

```
params have main grad=args.fp32 grad accumulation > 0,
            shard optimizer state=args.shard optimizer state > 0,
        )
    optimizer = smp.DistributedOptimizer(optimizer) # wrap the optimizer with smp.DistributedOptimizer
    lr scheduler = get learning rate scheduler(optimizer, args)
    if args.fp16:
        model.register post step hook(lambda model, optimizer:
optimizer.init master params())
    if args.enable memory profiling > 0:
        model.register post partition hook(
            lambda model, optimizer: memory status(msg="After partition")
        )
    # load after wrapping model and optimizer with smp Distributed...
    if args.load full or args.load partial:
        if args.load partial and args.load full:
            print(
                "Since both --load partial and --load_full set, will try
to load from full checkpoint."
                "If the intention is to load from partial checkpoint,
please don't set --load_full"
            )
        partial = not args.load full
        path = args.checkpoint dir if partial else args.model dir
        translate from hf = not partial
        model, optimizer, total steps, start train path index,
start_batch_index = load model and optimizer(
            path,
            model,
            optimizer,
            lr scheduler,
            partial,
            args,
            translate from hf=translate from hf,
            seq length=args.max context width,
            load model=True,
            load optimizer=args.load partial > 0,
            num params=num params,
        if args.save or verify ckptsum:
            filename = "saved sum" if args.load full else
"saved partial sum"
            load_and_verify_ckptsum(
                args, model, optimizer,
filename=os.path.join(args.model dir, filename)
    else:
        total steps = 0
        start train path index = 0
        start batch index = 0
```

```
start = time.time()
    total steps, throughput, loss = train(
        model,
        optimizer,
        lr scheduler,
        model config,
        start train path index,
        start batch index,
        num params,
        total steps,
        args,
    time to train = time.time() - start
    if args.ci:
        print(f"[SMP METRIC] GPT2 Time to train {time to train}")
        print(f"[SMP METRIC] GPT2 samples/second {throughput}")
       print(f"[SMP METRIC] GPT2_Loss__{loss}")
        if not args.load_partial and not args.load_full:
            assert time to train < args.time to train
            assert throughput > args.throughput
            if args.loss:
                assert loss < args.loss
    if args.save_final_full_model:
        # saves full model at the end
        base path = f"trained gpt nparams-{num params} steps-
{total steps}.pt"
        out path = os.path.join(args.model dir, base path)
        if args.save or verify ckptsum:
            # Save optimizer and model tensor sums and scalars before
saving
            save ckptsum(args, model, optimizer,
filename=os.path.join(args.model dir, "saved sum"))
        if smp.rdp rank() == 0:
            save(
                out path,
                model,
                optimizer,
                lr scheduler,
                model config,
                num params,
                total steps,
                -1,
                args,
                partial=False,
                translate to hf=smp.tp size() > 1,
                seq length=args.max context width,
            )
    smp.barrier()
    if smp.rank() == 0:
        print("SMP training finished successfully")
```

if __name__ == "__main__":