

Guidance to run Horovod TensorFlow ResNet50 model on Nvidia H100 GPUs

Log in to the ACES cluster and run the commands below.

```
$cd $SCRATCH
$mkdir horovod
$cd horovod
$git clone https://github.com/tensorflow/benchmarks
```

We used the Nvidia NGC TensorFlow container for this project. You can either download the container by yourself or use the container we stored in the shared directory.

Don't forget to set your directory in the job file where you saved the container.

#Option 1: Pull the container by yourself

```
$srun --time=02:00:00 --mem=100G --pty bash -i
$cd $SCRATCH
$mkdir container
$cd container
$export SINGULARITY_CACHEDIR=$TMPDIR
$module load WebProxy
$singularity pull docker://nvcr.io/nvidia/tensorflow:23.03-tf2-py3
```

#Option 2: Use the container we stored in a shared directory

the path to the container is /scratch/data/containers/tf2-23.03-py3.sif

```
$cd benchmarks
# create a slurm job file horovod-tf-h100-ngc.sh and copy and paste the content below to it.
$batch horovod-tf-h100-ngc.sh
```

```
#!/bin/bash

##NECESSARY JOB SPECIFICATIONS
#SBATCH --job-name=aces_h100
#SBATCH --time=01:00:00
#SBATCH --nodes=8
#SBATCH --ntasks-per-node=1
#SBATCH --cpus-per-task=8
#SBATCH --partition=gpu
#SBATCH --gres=gpu:h100:1
#SBATCH --mem=100G
#SBATCH --output=./result/aces_h100.%j
##SBATCH --nodelist=ac036,ac037,ac038,ac039,ac040,ac046,ac047,ac048
#SBATCH --exclusive
##SBATCH --reservation=workshop

echo $SLURM_JOB_ID

hostname

date

# load all the necessary modules
module purge
module load WebProxy
module load GCC/10.3.0 OpenMPI/4.1.1

cd /scratch/user/$USER/horovod/benchmarks

for i in {1..5}
do
    mpirun -np 8 -x NCCL_IB_DISABLE=0 -x NCCL_IB_CUDA_SUPPORT=1 -mca btl_tcp_if_include
ib0 -x NCCL_SOCKET_IFNAME=ib0 -x NCCL_DEBUG=INFO --bind-to none --map-by slot --mca
plm_rsh_args "-p 50000" singularity exec --nv
/scratch/data/containers/tf2-23.03-py3.sif python
scripts/tf_cnn_benchmarks/tf_cnn_benchmarks.py --variable_update=horovod
--model=resnet50 --batch_size=128
done

date
```