## 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.

```
#! /bin/bash
##NECESSARY JOB SPECIFICATIONS
#SBATCH --job-name=aces_h100
#SBATCH --time=01:00:00
```

```
#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 $SLURM JOB ID
hostname
date
module purge
module load WebProxy
module load GCC/10.3.0 OpenMPI/4.1.1
cd /scratch/user/$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
```