



Energy Efficiency and Performance Analysis of Monte-Carlo Event Generators for HEP

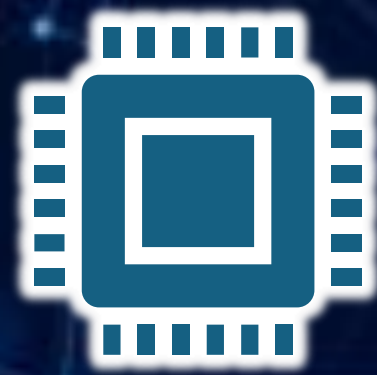
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OBJECTIVE



Study of power consumption and energy usage of **Madgraph** event generators

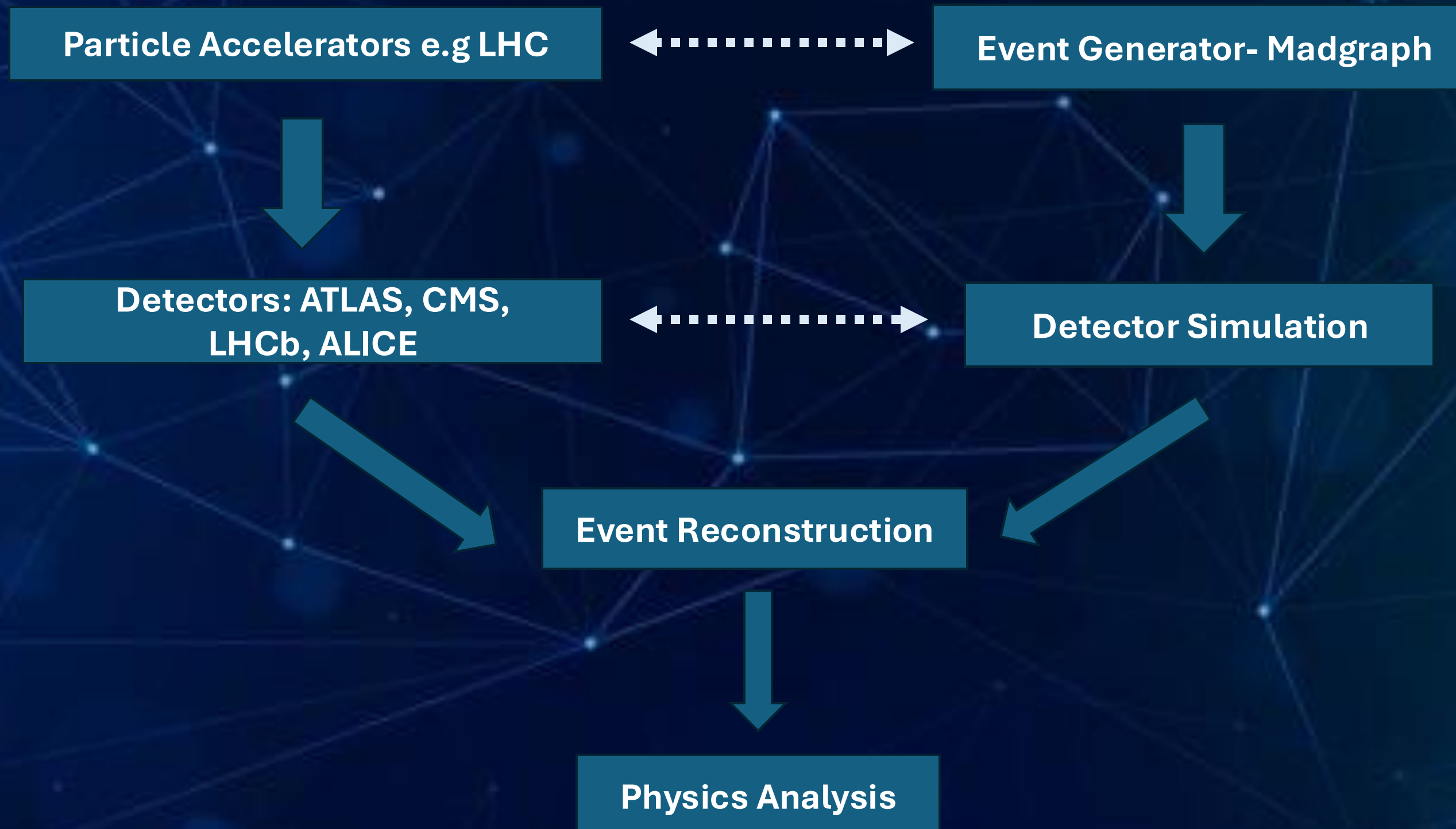


Using parallel **GPU** architecture



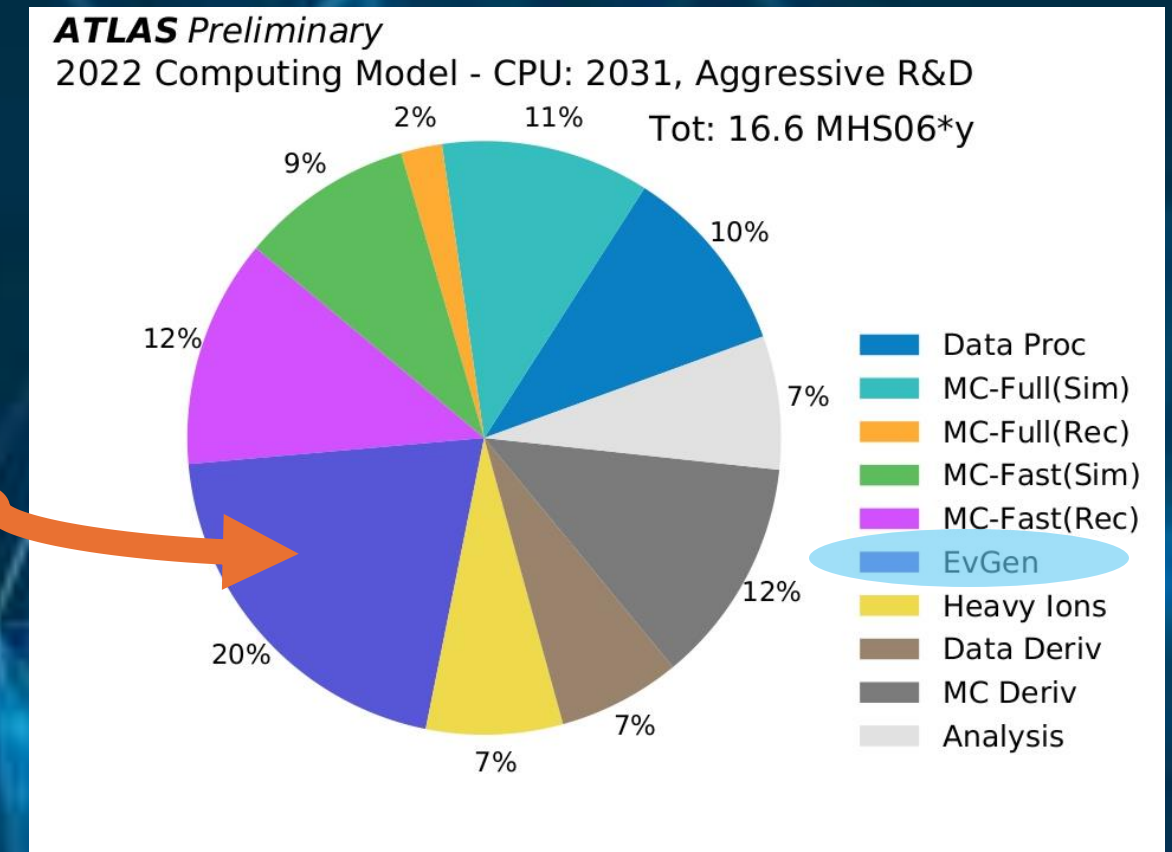
Vectorized CPU architecture like **AVX2** and **AVX512**

Big Picture



MOTIVATION

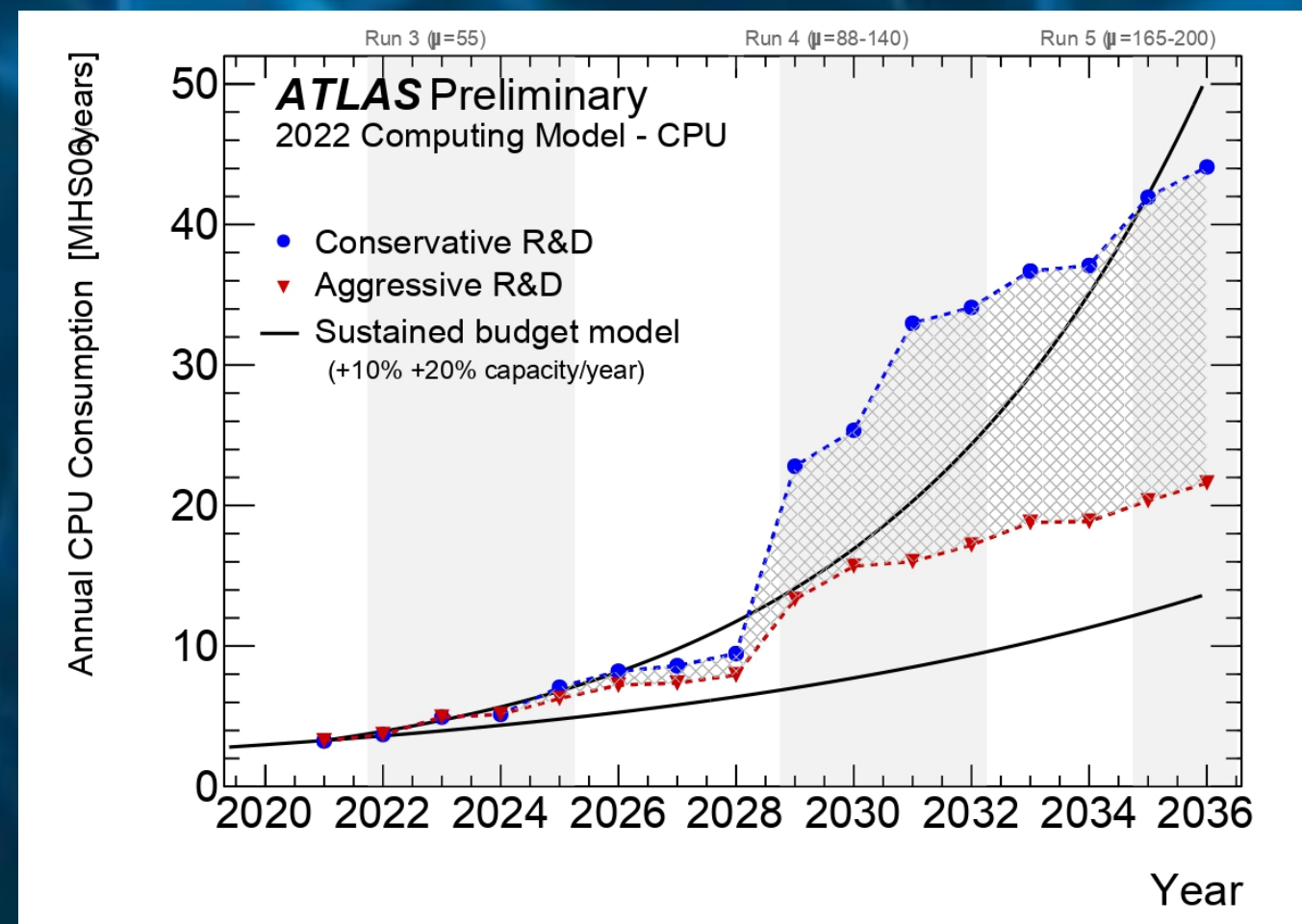
- Monte Carlo generators like MadGraph use 20% of computing resources
- More CPU or alternative hardware is needed
- Consider environmental impact



[ATLAS HL-LHC Computing Conceptual Design Report, 2022]

Approach

- Use SIMT GPUs like Nvidia and AMD
- Use SIMD CPUs like AVX2 and AVX-512



STEPS

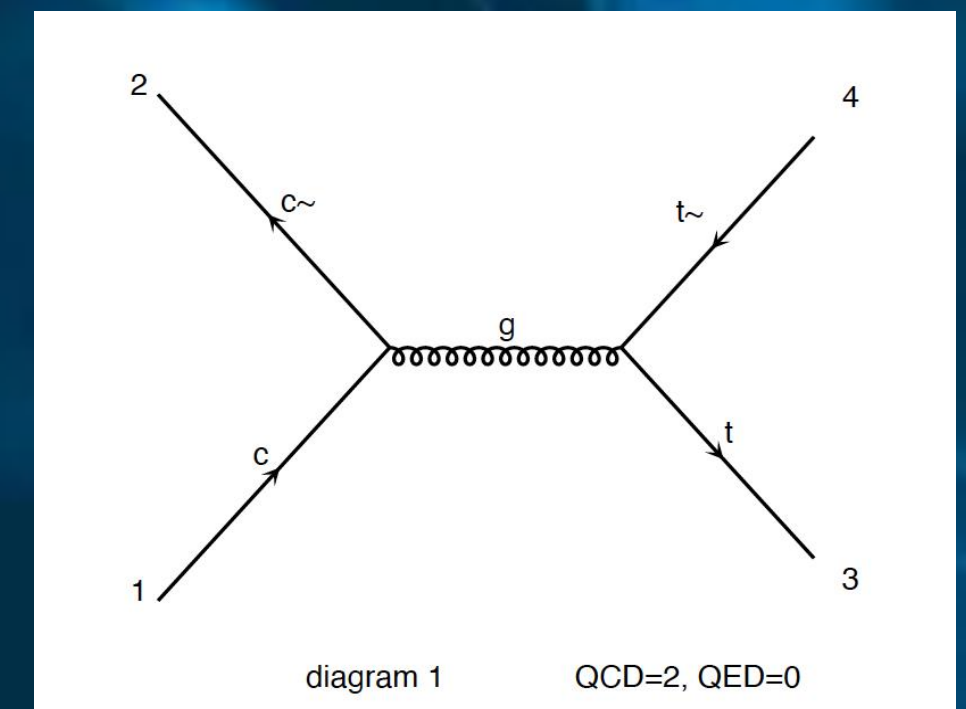
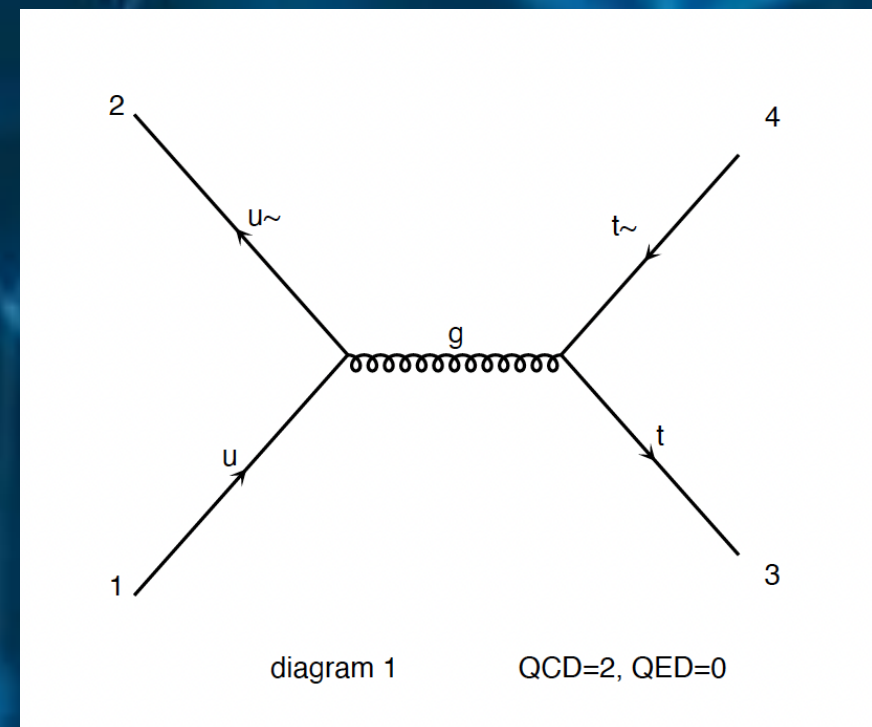
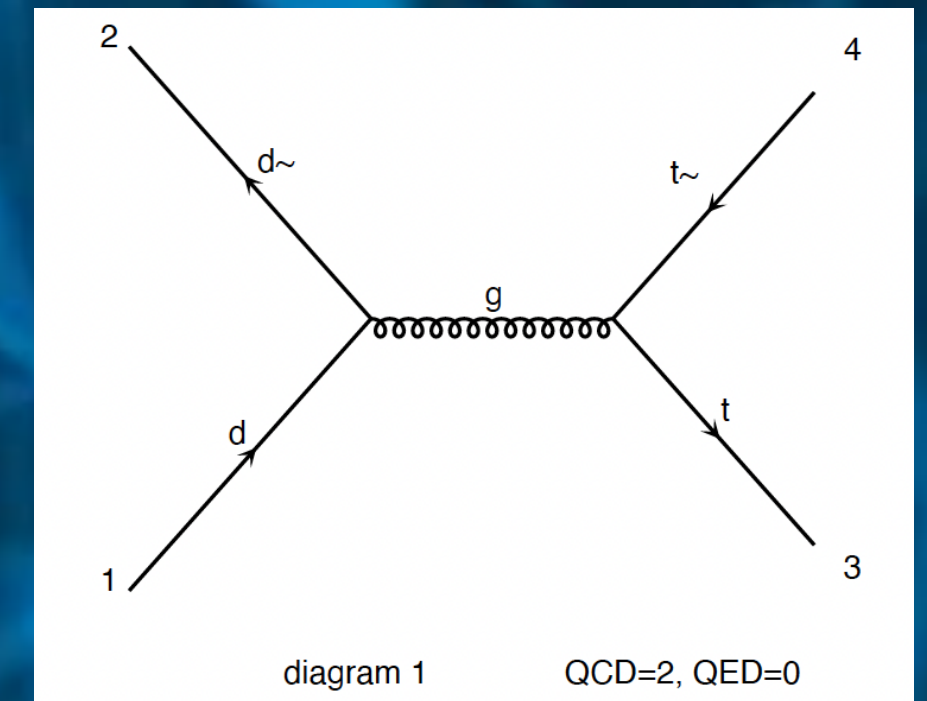
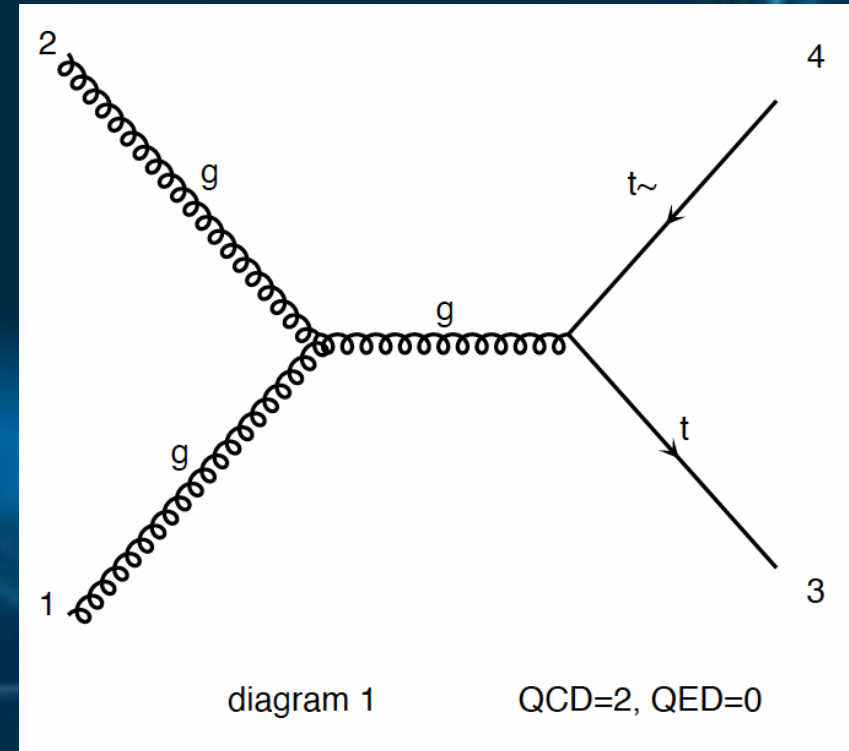
1. Single Thread Madevent:

- Madevent_gpu: to run processes on GPU
- Madevent_simd: to run processes on vector CPUs
- Runs the madevent generator for a single scattering sub-process

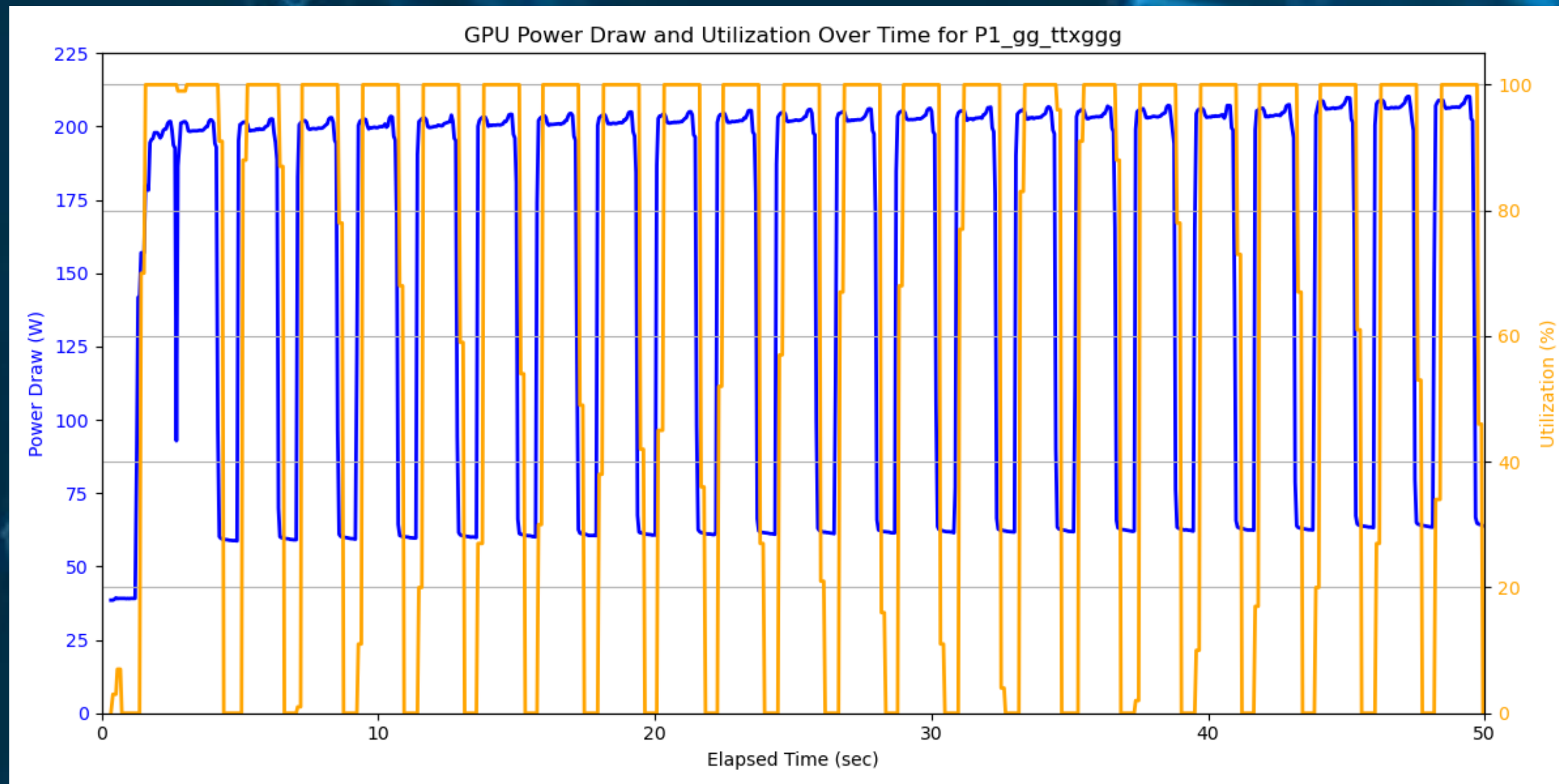
```
import model sm
generate p p > t t~
output madevent_gpu PROC_ppttx_gpu
launch PROC_pp_ttx_gpu
set cudacpp_backend cuda
```

2. WITH GRIDPACK: (Used by Experiments)

- Pre-compiled package
- Runs on multiple sub-processes



GPU Power and Utilization



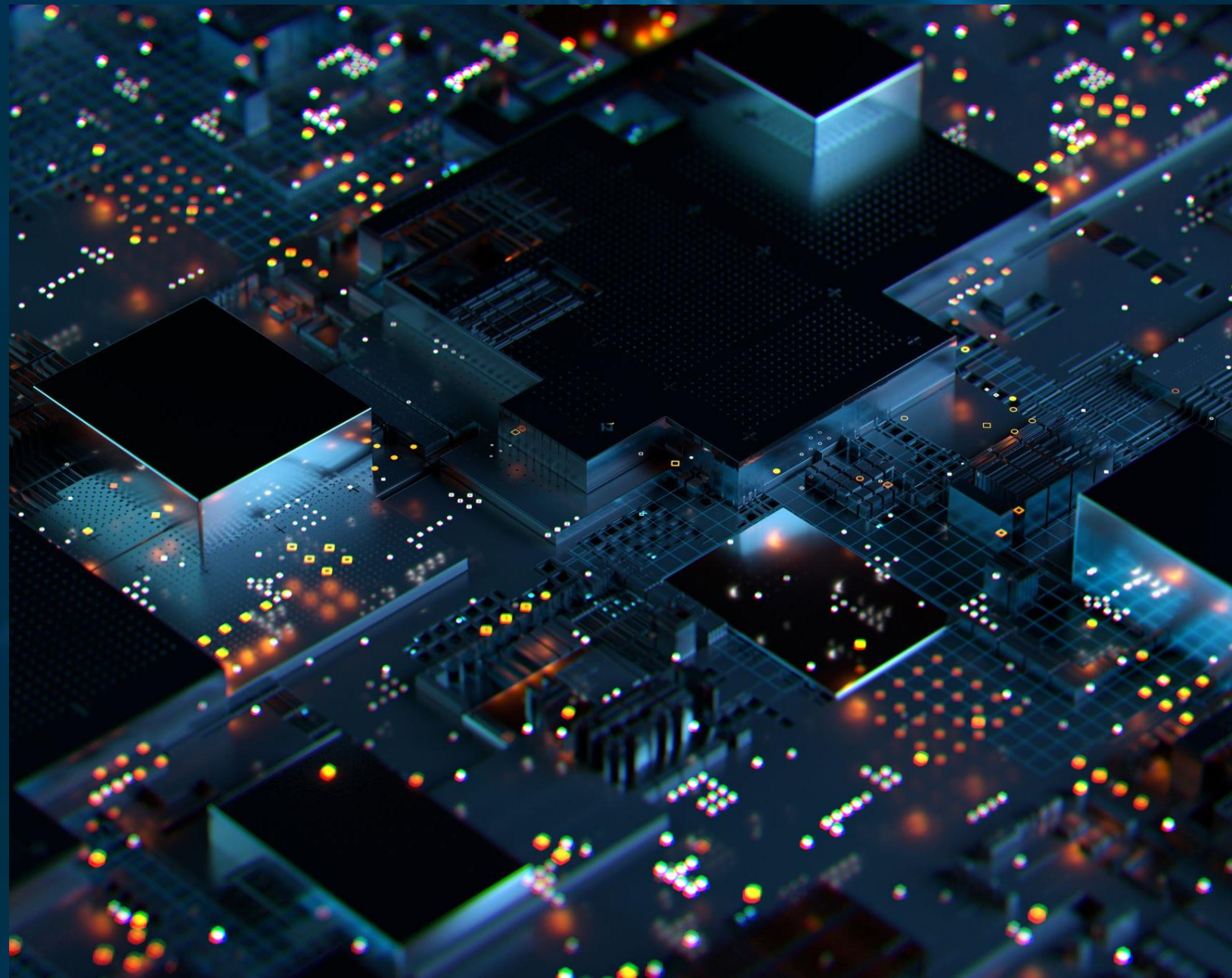
Single thread run done using Nvidia V100 GPU

Reliable Power Measurements on Gridpacks

- Use Cluster – CISM
- Execute commands via Slurm (with --exclusive flag)
- Generate gridpacks for various hardware and processes
 - Hardware: cuda, fortran, cppavx2
 - Node : mb-mil101
 - Processes: $g g > t t \sim ng$, $u u \sim > e^+ e^- ng$
- Developed automated scripts for gridpack generation and job launching
- Run the run.sh on a single node for:
 - Different values of threads and event number
 - Log start and end times of each execution



Future Steps



- Record the power consumption of GPU and CPU on gridpacks and produce plots
- Calculate energy by integrating power over time
- Calculate the CO2 emissions from the energy
<https://app.electricitymaps.com/map/72h/hourly>

Thank you!

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Scan Me!

Backups

Generating Gridpacks

```
genfile="generate_${process_name}_${backend}"

cat > "$genfile" <<EOF
generate $process
EOF

if [ $backend == "cuda" ]; then
|   echo "output madevent_gpu ${process_name}_${backend} -nojpeg" >> $genfile
else
|   echo "output madevent_simd ${process_name}_${backend} -nojpeg" >> $genfile
fi

echo "launch" >> $genfile
echo "set gridpack True" >> $genfile
echo "set cudacpp_backend $backend" >> $genfile

# Run MadGraph
../bin/mg5_aMC $genfile

mkdir -p gridpack_${backend}_${process_name}
mv ${process_name}_${backend}/run_01_gridpack.tar.gz gridpack_${backend}_${process_name}/
echo "Gridpack moved to gridpack_${backend}_${process_name}/"

echo "Ending full script at $timestamp"
```

Launching with Sbatch

```
for process in "${processes[@]"; do
  process_name=$(echo "$process" | xargs python -c "import sys; print(''.join(sys.argv[1:]).replace('>','_').replace('~','x').replace(' ','_'))")
  for backend in $backends; do
    echo "Submitting job for process: ${process_name} backend: $backend"
    if [ $backend == "cuda" ]; then
      sbatch --partition=gpu --exclusive --gres=gpu:1 --job-name=${process_name}_${backend} --output=logs/${process_name}_${backend}_%j.out producing_gridpack.sh $backend $process
    else
      sbatch --partition=gpu --job-name=${process_name}_${backend} --output=logs/${process_name}_${backend}_%j.out producing_gridpack.sh $backend $process
    fi
  done
done
```

```
# Define backends
backends="cuda fortran cppavx2"
processes=('g g > t t~' 'g g > t t~ g' 'g g > t t~ g g' 'u u~ > e+ e-' 'u u~ > e+ e- g' 'u u~ > e+ e- g g')
n_events=10000
sleep_time=200

nvidia-smi
cat /proc/cpuinfo

for process in "${processes[@]"; do
  process_name=$(echo "$process" | xargs python -c "import sys; print(''.join(sys.argv[1:]).replace('>','_').replace('~','x').replace(' ','_'))")
  for backend in $backends; do
    echo "=== Running for backend: $backend and process: $process_name ==="
    cd /home/ucl/cp3/rrichi/MG5_aMC_v3_6_3/work_dr/gridpack_${backend}_${process_name}/
    tar xzf run_01_gridpack.tar.gz
    for p in 1 2 4; do
      echo "Sleeping for $sleep_time s before running $backend with p: $p, for process: $process_name"
      sleep $sleep_time
      echo "Starting running with p: $p, for process: $process_name with $backend at $(date +%Y-%m-%d_%H-%M-%S)"
      ./run.sh -p $p $n_events 23
      echo "Ending running with p: $p, for process: $process_name with $backend at $(date +%Y-%m-%d_%H-%M-%S)"
    done
  done

  cd ..
done
```




CPU- Reliable?

No Because Shared!



Device I81887271383939

You can modify information and network configuration of the device.

i Information

Virtual machines 4

History

Name

Filter virtual machines

B9G47N3042

B9G47N3856

ITSCRD-V100

ITSCRD90

What is MadGraph?

Matrix Element Event Generator

Simulates Particle Collision

User defines initial and final particles

Generates Feynman Diagrams

Calculates physical quantities like cross sections

WHY MADGRAPH?



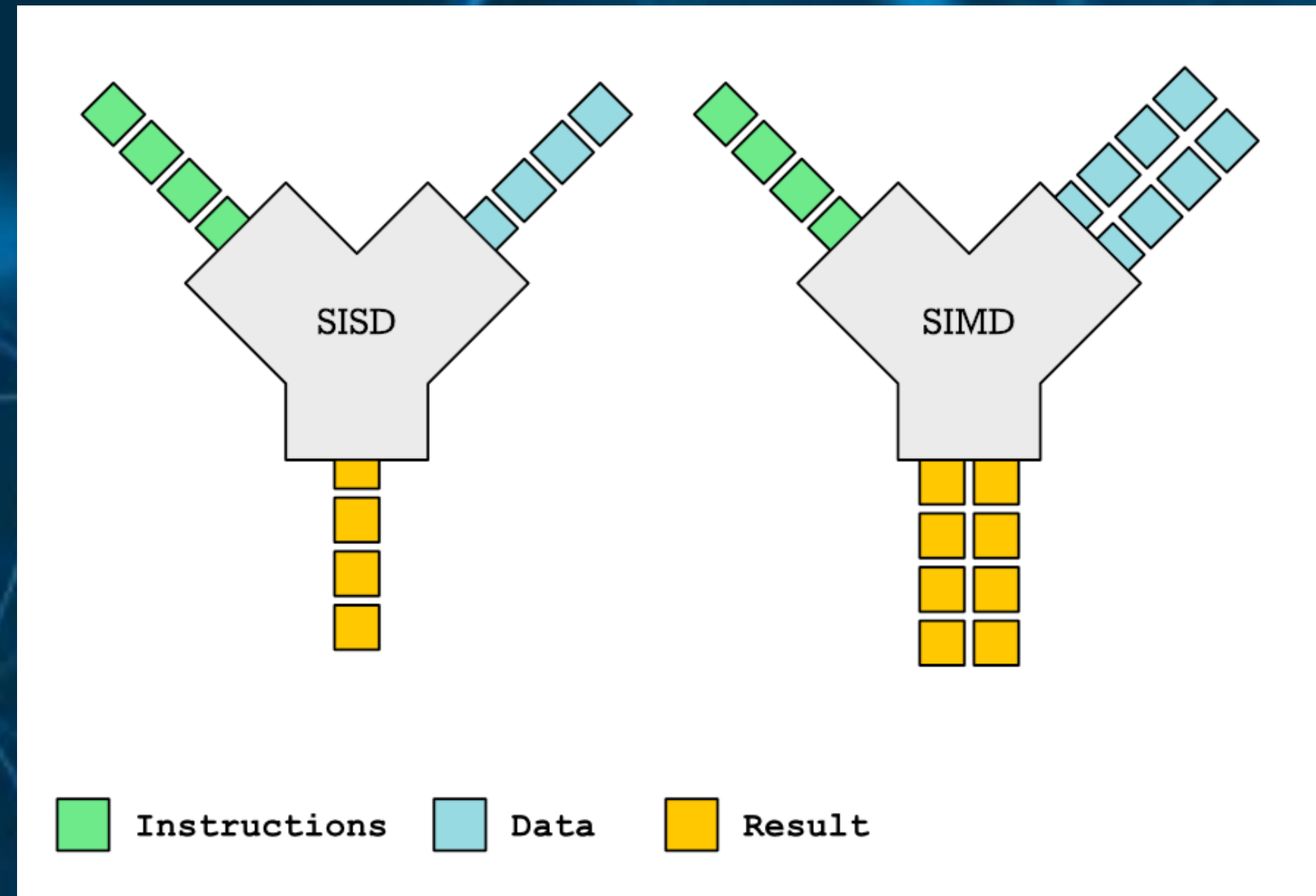
- **Vector CPUs & GPUs** are hard to utilize in HEP due to complex software workflows
- **MC detector simulations** involve **stochastic branching**, limiting data-parallel (SIMD) processing
- **Matrix Element Event Generators (like MadGraph)** compute scattering amplitudes at **independent phase-space points**. This makes them **ideal for SIMD & GPU acceleration**
- **Leverage hardware accelerators** to reduce compute time & energy consumption

SISD (Single Instruction Single Data):

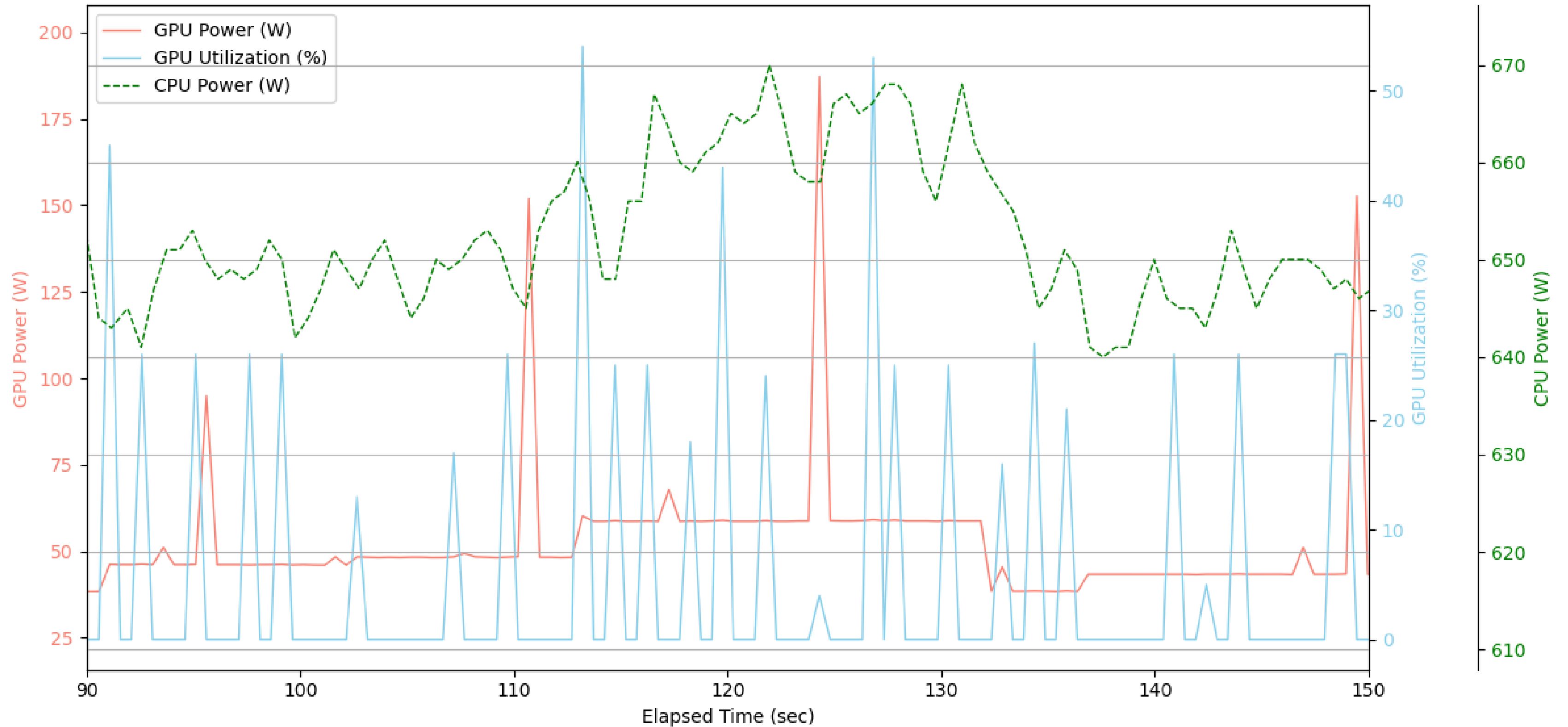
- Single processor executes one instruction at a time
- Requires more execution time
- Cannot scale well with increased computational needs

SIMD (Single Instruction Multiple Data):

- Process data points in parallel
- Excels in tasks requiring repetitive operations, reducing execution time
- Ensures proper memory alignment for faster execution



GPU(cuda) and CPU for ppttxgg



GPU(cuda) and CPU for ppttxggg

