# Sampled simulation with gem5



# What if the ROI is large

We now know how to skip the "unimportant" part of the simulation, but what if the important part of the simulation is too large?

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What if we are not facing this

What if we are not facing this

Kernel Boot Wider OS setup Benchmark application load ROI Finish and clean up
---

but actually facing this

# Sampling



# What is sampling?

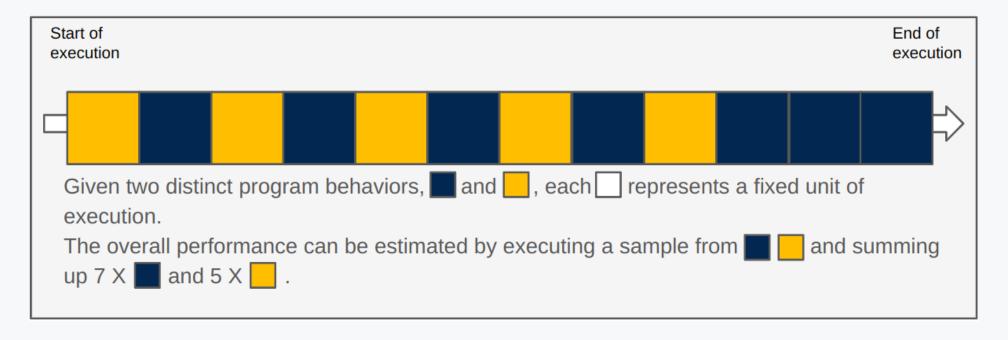
There are two major types of sampling:

- 1. Targeted sampling
- 2. Statistical sampling



# **Targeted Sampling**

Representative methodologies: SimPoint, LoopPoint



Targeted sampling selects samples based on specific characteristics that are discovered by analysis.

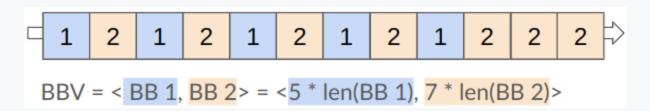


## **More About Targeted Sampling**

Well-known simulation sampling methods that use the targeted sampling approach include **SimPoint** and **LoopPoint**.

Both methods divide the whole program execution into regions that each execute a fixed number of instructions.

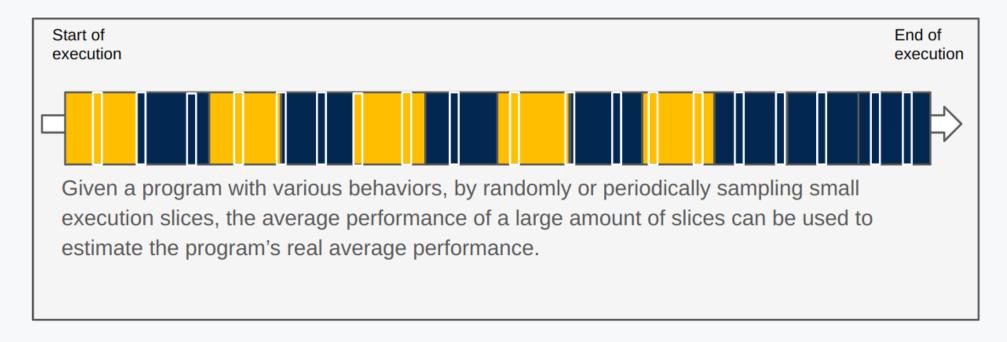
They use **basic block vectors**, which are recordings of the execution pattern of basic blocks within a region (period) of execution. A basic block vector can be used as the signature of the program behavior of the region. Here is an example of a basic block vector.



They use the basic block vectors to cluster and find the representative regions. They will predict the overall performance of the program by collecting only the representative regions' performance and sum them with weights.

## **Statistical Sampling**

Representative methodologies: SMARTS, FSA



Statistical sampling, as the name suggests, statistically selects its sampling units.



#### **More about Statistical Sampling**

The representative simulation sampling methods in statistical sampling include **SMARTS** and **FSA**.

Both methods periodically or randomly simulate in detail for a small amount of execution throughout the whole program execution and fast-forward between the detailed simulations.

They use the performance of the randomly distributed samples to predict the overall performance of the whole program execution.



#### What should we know before we apply the techniques

#### No matter how great a tool or a technique is, misusing it can be DANGEROUS

Before using any of the sampling techniques, we need to make sure the sampling technique works for our experiments.

For example, SimPoint is designed to work with single-threaded workloads only, so if our experiments require multi-threaded workloads, we should NOT use SimPoint with them.





#### What gem5 offers

In gem5, we have the infrastructure for

- 1. SimPoint (paper)
- 2. LoopPoint (paper)
- 3. ELFie (paper)
- 4. SMARTS (paper)
- 5. FSA (paper) (Might not be supported officially)



# **Targeted Sampling in gem5**



## Targeted Sampling in gem5

- gem5 provides infrastructure for **SimPoint** and **LoopPoint** to analyze the program, take checkpoints for the representative regions, and run the representative regions. Note that LoopPoint analysis support is not currently supported in gem5 v24.0 but is tested and prepared to be upstream in gem5 v24.1.
- gem5 also provides the infrastructure for **ELFies** to be executed in SE mode, but gem5 does not support creating ELFie files and weight information.



# SimPoint in gem5



#### **SimPoint**

As mentioned before, there are three steps to using SimPoint:

- 1. Analysis
- 2. Taking checkpoints
- 3. Run the regions

There are two key files related to using SimPoint in gem5:

- 1. <a href="mailto:src/python/gem5/utils/simpoint.py">src/python/gem5/utils/simpoint.py</a>
- 2. <a href="mailto:src/cpu/simple/probes/SimPoint.py">src/cpu/simple/probes/SimPoint.py</a>

We will be seeing them throughout this section.



## **SimPoint Analysis**

In gem5, we use the SimPoint probe listener object to collect the information SimPoint needs to cluster the regions.

This object is defined in <a href="mailto:src/cpu/simple/probes/SimPoint.py">src/cpu/simple/probes/SimPoint.py</a>.

The SimPoint probe listener has two parameters: interval and profile\_file.

- The [interval] takes a length as our definition of a region. It means that every time we execute n number of instructions, we see it as the end of a region. The default length is 100,000,000.
- The profile\_file takes a name for the output zip file. The default name is simpoint.bb.gz.

In order to use this probe listener object, we need to attach it to an ATOMIC CPU. It will start collecting information as soon as the simulation starts and will stop when the simulation ends. After exiting the simulation, there will be a zip file with the basic block vector information for each region under the simulation outdir directory.



#### **Hands-On Time!**

#### 01-simpoint

All materials can be found under <a href="materials/02-Using-gem5/09-sampling/01-simpoint">materials/02-Using-gem5/09-sampling/01-simpoint</a>. The completed version is under <a href="materials/02-Using-gem5/09-sampling/01-simpoint/complete">materials/02-Using-gem5/09-sampling/01-simpoint/simpoint/simpoint-simpoint/simpoint-simpoint-simpoint-simpoint/simpoint-simpoint/simpoint-simpoint/simpoint-s

#### Goal

- 1. Run the SimPoint analysis
- 2. Process the data to get the representative regions and their weights



Because the profiling and getting the baseline performance might take a while, we will run the simulation first with the following command.

```
gem5 -re --outdir=full-detailed-run-m5out full-detailed-run.py
gem5 -re --outdir=simpoint-analysis-m5out simpoint-analysis.py
```

In this exercise, we are trying to create SimPoints for a simple workload. The source code of the simple workload can be found in <a href="mailto:materials/02-Using-gem5/09-sampling/01-simpoint/workload/simple\_workload.c">materials/02-Using-gem5/09-sampling/01-simpoint/workload/simple\_workload.c</a>.

This simple workload allocates an array of one thousand 64 bit elements, assigns each a number, then sums it all up with one thousand iterations.

We can expect the program behavior of this workload to be really repetitive.



The script <u>materials/02-Using-gem5/09-sampling/01-simpoint/simpoint-analysis.py</u> uses the <u>SimPoint</u> probe listener object that we introduced earlier to collect basic block information for this simple workload.

The script <u>materials/02-Using-gem5/09-sampling/01-simpoint/simpoint-analysis.py</u> uses the <u>SimPoint</u> probe listener object that we introduced earlier to collect basic block information for this simple workload.

It connects the ATOMIC CPU core to the [SimPoint] probe listener using

```
processor.get_cores()[0].core.addSimPointProbe(1_000_000)
```

The definition of this <code>addSimPointProbe()</code> function can be found under <a href="mailto:src/cpu/simple/BaseAtomicSimpleCPU.py">src/cpu/simple/BaseAtomicSimpleCPU.py</a>.

In this example, we set the <u>interval\_length</u> as 1,000,000, which means that we define a region as 1,000,000 instruction executed (committed).



After the simulation finishes, we will see a zip file named simpoint.bb.gz under the simpoint-analysis-m5out folder.

We can unzip it with the following command:

```
gzip -d -k simpoint.bb.gz
```

After unzipping it, we can look at the [simpoint.bb] file.

This file contains all basic block vector information for the simple workload.

There are 9 regions found in this workload.

Each region has a basic block vector, which starts with a  $[\tau]$ .



As we can see, regions 2 to 9 have almost identical basic block vectors.

```
T:1900:222 :1901:222 :1902:999216 :1903:333
T:1900:222 :1901:222 :1902:999225 :1903:333
T:1900:222 :1901:222 :1902:999225 :1903:333
T:1900:222 :1901:222 :1902:999216 :1903:333
T:1900:222 :1901:222 :1902:999216 :1903:333
T:1900:222 :1901:222 :1902:999225 :1903:333
T:1900:222 :1901:222 :1902:999225 :1903:333
T:1900:222 :1901:222 :1902:999225 :1903:333
```

The similar basic block vectors indicate that program behavior for regions 2 to 9 are very similar. Therefore, we can expect to cluster regions 2 to 9 together and pick one region out of it to be our representative region. This region will represent the performance of all regions from region 2 to region 9.



Let's further understand what this line means.

T:1900:222 :1901:222 :1902:999216 :1903:333

- T means the start of the region's basic block vector.
- :1900:222 means that basic block 1900 executed (committed) 222 instructions. The key here is that 222 is NOT the number of times the basic block has been executed, but the number of times the basic block has been executed multiplied by the total instructions in the basic block . If we sum up all the instructions that have been executed by the basic blocks, we will get roughly the length of the region. 222 + 222 + 999216 + 333 = 999993.

The next step is to use this information to cluster the regions and find the representative regions.

There are many methods to do it. In this exercise, we will be using the SimPoint3.2 tool that was provided by the SimPoint paper's authors. <u>Link</u> to the tool.



The tool is already compiled under <u>materials/02-Using-gem5/09-sampling/01-simpoint/simpoint</u>. We also provided a runscript with the following command in <u>materials/02-Using-gem5/09-sampling/01-simpoint/simpoint3.2-cmd.sh</u>.

```
/workspaces/2024/materials/02-Using-gem5/09-sampling/01-simpoint/simpoint \
-inputVectorsGzipped -loadFVFile simpoint-analysis-m5out/simpoint.bb.gz -k 5 -saveSimpoints \
results.simpts -saveSimpointWeights results.weights
```

Let's look at this command.

It passes in simpoint.bb.gz to the tool.

It sets the number of clusters expected to 5 using [-k 5].

It saves the SimPoint information in results.simpts and their weights in results.weight.



After running the command with [./simpoint3.2-cmd.sh], we will see the following.

```
#results.simpts # results.weights
2 0 0.666667 0
1 1 0.222222 1
0 3 0.11111 3
```

This means that it found 3 SimPoints out of this program.

Region 2 is SimPoint 0, which has a weight of 0.666667.

Region 1 is SimPoint 1, which has a weight of 0.222222.

Region 0 is SimPoint 3, which has a weight of 0.111111.

There are not 5 clusters because the algorithm found that 3 clusters are enough to represent all the program behaviors.

The SimPoint tag number might not be continuous because it is the tag number for the cluster.



# **SimPoint Checkpoint**

Now that we have the representative regions and their weights, we will need to find a way to get to those SimPoints.

As introduced in <u>08-accelerating-simulation</u>, there are two methods to get to the region of interest. With SimPoint, we usually use checkpointing.

To do it in gem5, we will be using <u>set\_se\_simpoint\_workoad</u>, the <u>SimPoint class</u>, and the <u>simpoints\_save\_checkpoint\_generator</u>.



#### **Hands-On Time!**

## 01-simpoint

All materials can be found under <u>materials/02-Using-gem5/09-sampling/01-simpoint</u>. The completed version is under <u>materials/02-Using-gem5/09-sampling/01-simpoint/complete</u>. We will not be modifying any scripts.

#### Goal

• Use the output file from the SimPoint3.2 tool to create checkpoints for all 3 SimPoints

Because it might take some time to take the checkpoints, let's start taking checkpoints with the following command:

gem5 -re --outdir=simpoint-checkpoint-m5out simpoint-checkpoint.py



Let's look at <u>materials/02-Using-gem5/09-sampling/01-simpoint/simpoint-checkpoint.py</u>.

#### There are three key parts:

```
# key 1: pass in the representative regions information
from gem5.utils.simpoint import SimPoint
simpoint_info = SimPoint(
    simpoint_interval=1_000_000,
    simpoint_file_path=Path("/workspaces/2024/materials/02-Using-gem5/09-sampling/01-simpoint/results.simpts"),
    weight_file_path=Path("/workspaces/2024/materials/02-Using-gem5/09-sampling/01-simpoint/results.weights"),
    warmup_interval=1_000_000
# key 2: pss in the SimPoint object to set up the workload
board.set_se_simpoint_workload(
    binary=BinaryResource(local_path=Path("/workspaces/2024/materials/02-Using-gem5/09-sampling/01-simpoint/workload/simple_workload").as_posix()),
    simpoint=simpoint_info
# key 3: register exit event handler to take the checkpoints
simulator = Simulator(
    board=board,
    on_exit_event={
        ExitEvent.SIMPOINT_BEGIN: simpoints_save_checkpoint_generator(dir, simpoint_info)
    },
```



Let's zoom in to key 1.

```
from gem5.utils.simpoint import SimPoint
simpoint_info = SimPoint(
    simpoint_interval=1_000_000,
    simpoint_file_path=Path("/workspaces/2024/materials/02-Using-gem5/09-sampling/01-simpoint/results.simpts"),
    weight_file_path=Path("/workspaces/2024/materials/02-Using-gem5/09-sampling/01-simpoint/results.weights"),
    warmup_interval=1_000_000
)
```

The name of the parameters for <code>simpoint\_interval</code>, <code>simpoint\_file\_path</code>, and <code>weight\_file\_path</code> should speak for themselves, but <code>warmup\_interval</code> is new to us in this section.

As mentioned in <u>08-accelerating-simulation</u>, we should expect most of the micro-architectural states to be cold when restoring a checkpoint. Therefore, we need to reserve a warm up period for our region of interest to warm up the micro-architectural states of the system.

As such, we need to set up a warmup\_interval value here, so when we restore the checkpoint, we can expect to have this length of simulation to warm up our micro-architectural states.

The SimPoint class can also detect how much space we have for the warm up period. If there is not enough space, it will automatically shrink the warm up period of that SimPoint to the space available. For example, if we have a SimPoint that starts at the beginning of the program, then the warm up period for it should be 0 instructions.

The SimPoint class also automatically sorts the SimPoints based on when they start in terms of instructions.

It also provides getter functions that might be helpful, such as  $[get_weight_list()]$  that returns the list of the weights for each SimPoint.

More information can be found in the **SimPoint** class definition.



Let's just trust keys 2 and 3 because they both rely on the information passed in from the SimPoint object.

We can declare where we want to store the SimPoint checkpoints by passing in the destination path to the simpoints\_save\_checkpoint\_generator as a parameter. For more details, we can find the generator in <a href="mailto:src/python/gem5/simulate/exit">src/python/gem5/simulate/exit</a> event <a href="mailto:generators.py">generators.py</a>.



After we finish checkpointing, we should find all of the SimPoint checkpoints under <a href="materials/02-">materials/02-</a>
<a href="materials/02-">Using-gem5/09-sampling/01-simpoint/simpoint-checkpoint</a> because we pass in <a href="materials/02-">simpoint-checkpoint</a> because we pass in <a href="materials/02-">simpoint-checkpoint</a> to the <a href="materials/02-">simpoints\_save\_checkpoint</a> as our directory to save the SimPoint checkpoints.

There should be three checkpoint folders called <code>cpt.SimPoint0</code>, <code>cpt.SimPoint1</code>, and <code>cpt.SimPoint2</code>.

We can now use them to run our SimPoints for the simple workload.



## **Running SimPoint**

Now we have the representative SimPoint checkpoint, we can run them with the systems we want to measure and predict the overall performance of running the whole simple workload.

SimPoint relies on the weights we got from the analysis stage to do the prediction. The weight is calculated with

Weight of cluster 
$$i = \frac{\text{Number of regions in cluster } i}{\text{Total number of regions in the workload execution}}$$

For example, if we want to calculate the predicted IPC, we should use

$$ext{Predicted IPC} = \sum_{i=1}^n ( ext{Weight of cluster } i imes ext{IPC of cluster } i)$$



All materials can be found under <u>materials/02-Using-gem5/09-sampling/01-simpoint</u>. The completed version is under <u>materials/02-Using-gem5/09-sampling/01-simpoint/complete</u>. We will still not modify any scripts.

Unlike the simple systems we used for SimPoint analysis and SimPoint checkpointing, we now need to use the systems for which we actually want to measure the performance.

We will be using <u>materials/02-Using-gem5/09-sampling/01-simpoint/simpoint-run.py</u> to run our SimPoints.

For our baseline, we are using <u>materials/02-Using-gem5/09-sampling/01-simpoint/full-detailed-run.py</u>, which runs the whole simple workload with the detailed system.

Let's start by running the SimPoints before explaining how it works due to time constraints. We provided a runscript to run all three in <a href="materials/02-Using-gem5/09-sampling/01-simpoint/run-all-simpoint.sh">materials/02-Using-gem5/09-sampling/01-simpoint/run-all-simpoint.sh</a>

./run-all-simpoint.sh



Let's look at <u>materials/02-Using-gem5/09-sampling/01-simpoint/simpoint-run.py</u>. It has a detailed system that matches the one that is used in our baseline <u>materials/02-Using-gem5/09-sampling/01-simpoint/full-detailed-run.py</u>.

There are a few key points that we want to look at. Let's start with the ones we are familiar with.

```
# key 1:
from gem5.utils.simpoint import SimPoint
simpoint_info = SimPoint(
    simpoint_interval=1_000_000,
    simpoint_file_path=Path("/workspaces/2024/materials/02-Using-gem5/09-sampling/01-simpoint/results.simpts"),
    weight_file_path=Path("/workspaces/2024/materials/02-Using-gem5/09-sampling/01-simpoint/results.weights"),
    warmup_interval=1_000_000
)
# key 2:
board.set_se_simpoint_workload(
    binary=BinaryResource(local_path=Path("/workspaces/2024/materials/02-Using-gem5/09-sampling/01-simpoint/workload/simple_workload").as_posix()),
    simpoint=simpoint_info,
    checkpoint=Path(f"simpoint-checkpoint/cpt.SimPoint{args.sid}")
)
```

Here, we pass in the SimPoint information and use it to set up a workload that restores the checkpoint with our target SimPoint id.

```
# kev 3:
def max inst():
    warmed_up = False
    while True:
        if warmed up:
            print("end of SimPoint interval")
            vield True
        else:
            print("end of warmup, starting to simulate SimPoint")
            # Schedule a MAX_INSTS exit event during the simulation
            simulator.schedule_max_insts(
                board.get_simpoint().get_simpoint_interval()
            m5.stats.dump()
            m5.stats.reset()
            yield False
simulator = Simulator(
    board=board.
    on_exit_event={ExitEvent.MAX_INSTS: max_inst()},
```

Here is an exit event handler we defined for handling the warm up period and detailed simulation period.

It schedules an end for the SimPoint after the warm up period. It also dumps and resets the stats for detailed measurement.



```
# key 4:
warmup_interval = board.get_simpoint().get_warmup_list()[args.sid]
if warmup_interval == 0:
    warmup_interval = 1
print(f"Starting Simulation with warmup interval {warmup_interval}")
simulator.schedule_max_insts(warmup_interval)
simulator.run()

print("Simulation Done")
print(f"Ran SimPoint {args.sid} with weight {board.get_simpoint().get_weight_list()[args.sid]}")
```

Before the starting the simulation, we need to set up an exit event to indicates the end of the warm up period.

We use the simulator.schedule\_max\_insts() function to do so.

We can use [get\_warmup\_list()] from the [SimPoint] object to get the warm up period length for each SimPoint.

Here is one limitation of the  $[simulator.schedule_max_insts()]$  function. If it gets the value [0], there won't be any exit event scheduled, so if the warm up period length is [0], we have to set it to [1].



After setting up the above keys, the script is ready to run the SimPoint. Note that each simulation can only run one SimPoint.

After the running the SimPoints, we should see the output folder [simpoint0-run], [simpoint1-run], and [simpoint2-run].

Also, we should have the baseline output in [full-detailed-run-m5out].

Let's try to use the SimPoints' performance to predict the overall IPC!

We have a Python script ready to do the prediction in <u>materials/02-Using-gem5/09-sampling/01-simpoint/predict\_overall\_ipc.py</u>.

We can run it with

python3 predict\_overall\_ipc.py



## 01-simpoint

We should see something like this

```
predicted IPC: 1.2577933618669999
```

actual IPC: 1.247741

relative error: 0.8056449108428648%

The Python script reads the IPC from our baseline.

It also reads the detailed simulation period's IPC from all our SimPoints' stats files.

Then it performs the calculation with

```
predicted_ipc = 0.0
for i in range(num_simpoints):
    predicted_ipc += simpoint_ipcs[i] * simpoint_weights[i]
```

As the output suggests, the relative error between the predicted IPC and the actual baseline IPC is around 0.81%.

## **Summary in SimPoint**

Congratulations! We walked through the whole process of sampling with the SimPoint method.

Let's recap what we did:

- 1. Analyzing the program
- 2. Checkpointing the representative regions
- 3. Running the regions and predicting the performance

Good news: this process is very similar for the majority of the methods in targeted sampling. Therefore, if we know how to do sampling with the SimPoint method, it should not be hard for us to use other methods, such as LoopPoint, which enable multi-threaded workload sampling.



# **LoopPoint and ELFies in gem5**



## **LoopPoint and ELFies**

LoopPoint is similar to SimPoint with some key differences.

LoopPoint uses the number of times a loop is executed to mark the regions instead of using instructions executed.

Therefore, we need to collect the loop execution information in the analysis stage in addition to the basic block execution information.

Other than this, it is very similar to SimPoint in terms of the process (the 3 step process we did in 01-simpoint).



## **LoopPoint and ELFies**

ELFies is a checkpointing method that creates checkpoint executables out of a large workload execution. It can be used with LoopPoint to create executables of the representative regions.

Like LoopPoint, it uses the number of times a loop has been executed to mark the beginning and the end of the region of interest, so we need that information to execute ELFies in gem5.

<u>Link</u> for more information about LoopPoint and ELFies.



#### **ELFies**

gem5 does not produce ELFies, but we have support to run ELFies in SE mode. All the weights and loop information should come with the ELFies workload.

We can run ELFies with the **ELFielnfo class**.

There is an example in <u>materials/02-Using-gem5/09-sampling/02-elfies/run-elfies.py</u>

We can run it with the following command but it is not suggested because it will take too long.

gem5 -re run-elfies.py

It is a 8-threaded experiment with a detailed system that might run for eight hundred million instructions so it will take some time to finish.

We have a completed m5out under the <u>materials/02-Using-gem5/09-sampling/02-elfies/complete/m5out</u> if you are interested in looking at the output.



## **ELFies Example**

We provide some ELFie workloads, such as <u>wrf-s.1\_globalr13</u>, in gem5 resources. Its marker (loop execution) information can be found in <u>elfie-info-wrf-s.1\_globalr13</u>.

We can set up the ELFies workload as a normal executable.

```
board.set_se_binary_workload(
    binary=obtain_resource("wrf-s.1_globalr13")
)
```

Then, we can set up the beginning and end marker with

```
from gem5.resources.elfie import ELFieInfo
elfie = ELFieInfo(start = PcCountPair(int("0x100b643",0),1), end = PcCountPair(int("0x526730",0),297879) )
elfie.setup_processor(
    processor = processor
)
```



## **ELFies Example**

After getting to the beginning and end markers, the simulator will raise a SIMPOINT\_BEGIN exit event. However, there is no default exit event handler for this, so we will need to define our own.

```
def start_end_handler():
    # This is a generator to handle exit event produced by the
    # start marker and end marker.
    # When we reach the start marker, we reset the stats and
    # continue the simulation.
    print(f"reached {targets[0]}\n")
    print("now reset stats\n")
    m5.stats.reset()
    print("fall back to simulation\n")
    yield False
    # When we reach the end marker, we dump the stats to the stats file
    # and exit the simulation.
    print(f"reached {targets[1]}\n")
    print("now dump stats and exit simulation\n")
    m5.stats.dump()
    yield True
simulator = Simulator(
    board=board,
    on_exit_event={
        ExitEvent.SIMPOINT_BEGIN : start_end_handler()
```



## **Summary of LoopPoint and ELFies**

After we finish running all of the ELFies for a macro-benchmark, we can use the weights provided by the ELFie files to predict the overall performance like we did for the SimPoint example.

Now we covered all the targeted sampling methods that are supported in gem5, let's dive into statistical sampling!



## Statistical Sampling in gem5



SMARTS is one of the statical sampling methods.

It uses a statistical model to predict the overall performance with randomly or periodically selected samples.

We use small samples distributed throughout the whole program execution to predict the average performance. We can also use the average performance to predict the overall performance, such as runtime.

Before running the simulation, we need to decide a few statistical parameters.

- n: the number of the samples. It is a number of count.
- k: the systematic sampling interval. It is a number of count.
- [U]: the sampling unit size. It is a number of instructions to execute.
- $\bullet$  [w]: the length of the detailed warm up period. It is a number of instructions to execute.



### **Hands-On Time**

#### 03-SMARTS

All materials can be found under <u>materials/02-Using-gem5/09-sampling/03-SMARTS</u>. The completed version is under <u>materials/02-Using-gem5/09-sampling/03-SMARTS/complete</u>. We will not modify any scripts.

<u>materials/02-Using-gem5/09-sampling/03-SMARTS/SMARTS.py</u> is an example of how to use exit event handler to perform SMARTS.

We can run it with the following command

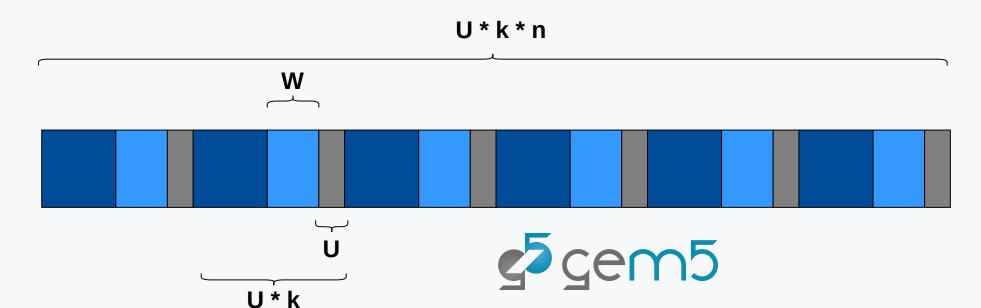
gem5 -re SMARTS.py

This script will use SMARTS on the workload from 01-simpoint, so we can use the baseline performance from <a href="mailto:materials/02-Using-gem5/09-sampling/01-simpoint/full-detailed-run.py">materials/02-Using-gem5/09-sampling/01-simpoint/full-detailed-run.py</a> to validate our predicted performance using SMARTS.



Let's look at the smarts\_generator with the statistical parameters.

- n: the number of the samples. It is a number of count.
- [k]: the systematic sampling interval. It is a number of count.
- [U]: the sampling unit size. It is a number of instructions executed.
- W: the length of the detailed warm up period. It is a number of instructions executed.



In order to determine  $\[mu]$  and  $\[mu]$ , we first need to determine our ideal  $\[mu]$ . In the SMARTS paper, they define a large sample size as n>30. For this exercise, let's set our ideal  $\[mu]$  as 50.

After setting up the ideal [n], we can use it with the instructions executed in the workload to determine the ideal [k].

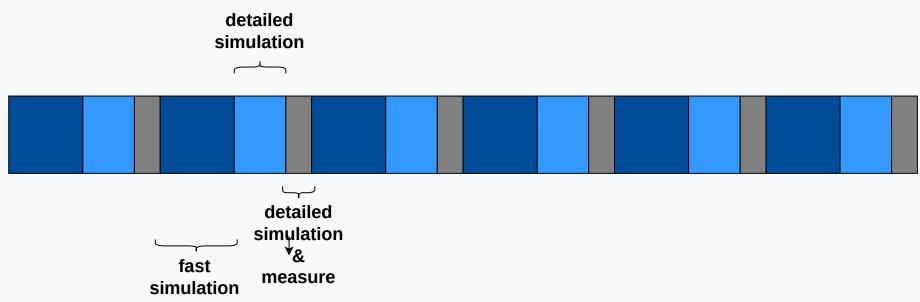
Here is an example of how it can be done in Python.

```
program_length = 9115640
ideal_region_length = math.ceil(program_length/50)
ideal_U = 1000
ideal_k = math.ceil(ideal_region_length/ideal_U)
ideal_W = 2 * ideal_U
```

With these parameters, we will have about 50 samples. Each sample has a gap (k-1)\*U=182000 instructions from each other. Each sample has 1000 instructions in detailed measurement and 2000 instructions of warm up period.

This SMARTS exit generator only works with SwitchableProcessor.

When it reaches to the start of the detailed warmup part, it resets the stats; then it switches the core type and schedule for the end of the warmup part and the end of the interval. When it reaches to the end of the detailed warmup part, it resets the stats. When it reaches to the end of the detailed simulation, it dumps the stats; then it switches the core type and schedule for the start of the next detailed warmup part.





Now we understand what the <code>smarts\_generator</code> will do, let's look at the system we are using in <a href="materials/02-Using-gem5/09-sampling/03-SMARTS/SMARTS.py">materials/02-Using-gem5/09-sampling/03-SMARTS/SMARTS.py</a>.

It uses exactly the same system as <u>materials/02-Using-gem5/09-sampling/01-simpoint/full-detailed-run.py</u>, except that it uses a <u>SimpleSwitchableProcessor</u> to switch between ATOMIC CPU and O3 CPU for fast-forwarding and detailed simulation.

If the simulation is run with

gem5 -re SMARTS.py

we can run <u>materials/02-Using-gem5/09-sampling/03-SMARTS/predict\_ipc.py</u> to predict the overall IPC and calculate the relative error with the baseline IPC.

python3 predict\_ipc.py



What <u>materials/02-Using-gem5/09-sampling/03-SMARTS/predict\_ipc.py</u> does is read the IPC from each sample, sum the IPCs, and divide by the total number of samples to calculate the average IPC of the program.

Average IPC 
$$=rac{\sum_{i=1}^{n} ext{IPC}_{i}}{n}$$

Here is what we expect to see

Number of samples: 50

Predicted Overall IPC: 1.2563117400000001

Actual Overall IPC: 1.247741

Relative Error: 0.6869005667041583%

As the output suggests, the relative error between the predicted IPC and the actual baseline IPC is around 0.69%.



## **Summary**

Now we experimented both targeted and statistical sampling in gem5, let's end it with the trade-offs.

Trade-off Method	Speedup	Resource Requirement	Prediction accuracy	Cost to change configuration	Ability to compare samples
Targeted Sampling	$\frac{S_{d/f}}{(S_{d/f} - 1) \times (n/N) + 2}$ $\frac{N}{n}$	Storage	It is impacted by the cluster signature and clustering algorithm	Almost no cost to change hardware configurations, but big cost to change software configurations	Allows directly comparison between samples
Statistical Sampling	$\frac{S_{d/f}}{(S_{d/f}-1)\times (n/N)+1}$	KVM & parallel resources	It is impacted by the input parameters	Equal cost to change both hardware and software configurations	Do not have comparable samples



## **Question in Mind**

When we run the SimPoints, we need to run a separate simulation for each SimPoint. What if we could do it all in one simulation?

