

HPC Community Tools

Presenter: Giacomo Capodaglio Oct 21-23, 2025 AMD @ Tsukuba University



HPC Community Tools

- TAU
 - https://www.cs.uoregon.edu/research/tau/home.php
- HPCToolkit
 - https://hpctoolkit.gitlab.io/
- Score-P
 - https://www.vi-hps.org/projects/score-p/overview/overview.html
- Likwid
 - https://docs.hpc.gwdg.de/software_stacks/applications/likwid/index.html
- Scalasca
 - https://scalasca.org/
- Grafana
 - https://grafana.com/
- Perfetto
 - https://ui.perfetto.dev/

TAU

TAU

- Tuning and Analysis Utilities, developed at University of Oregon
- Scalable and flexible performance analysis toolkit
- Automatic instrumentation through Program Database Toolkit (PDT) for routines, loops, I/O, memory, phases, etc.

Installing TAU

Installed from the public mirror on <u>GitHub</u>

```
(currently using GIT COMMIT="23a56e2a1a728e99ff03341c30f9d24892c5952b"):
    git clone https://github.com/UO-OACISS/tau2.git
    cd tau2
    git checkout $GIT COMMIT
    wget http://tau.uoregon.edu/ext.tgz
    tar zxf ext.tgz
# in the latest version of TAU -roctrace and -rocprofiler are mutually exclusive and -roprofsdk should be used
ROCM FLAGS="-rocm=${ROCM PATH} -hip=${ROCM PATH} -rocmsmi=${ROCM PATH} -roctracer=${ROCM PATH} -
rocprofiler=${ROCM PATH}"
      result=`echo $ROCM VERSION | awk '$1>6.1.2'` && echo $result
      if [[ "${result}" ]]; then # ROCM VERSION >= 6.2
         ROCM FLAGS="-rocm=${ROCM PATH} -hip=${ROCM PATH} -rocmsmi=${ROCM PATH} -rocprofsdk=${ROCM PATH}"
      fi
                                                               TAU creates a library for every configuration
    # configure with: MPI OMPT OPENMP PDT ROCM
    ./configure -c++=amdclang++ -fortran=gfortran -cc=amdclang -prefix=${TAU_PATH} -zlib=download -otf=download
                 -unwind=download -bfd=download ${ROCM FLAGS} -mpi -ompt -openmp -pdt=${PDT PATH} -iowrapper
```

How to use TAU

- There are mainly two approaches to use an application with TAU and GPUs
 - Use TAU Wrappers
 - For C: replace the compiler with tau cc.sh
 - For C++: replace the compiler with tau_cxx.sh
 - For Fortran: replace the compiler with tau_f90.sh
 - Dynamic instrumentation, for example:
 - srun -n 2 -gpus 2 tau_exec –T rocm,rocprofiler -rocm ./test

TAU example in HPCTrainingExample repo

```
cd ${REPO_DIR}/HIP/jacobi
module load rocm
module load openmpi
module load tau
export TAU_PROFILE=${TAU_PROFILE}
export TAU_TRACE=${TAU_TRACE}
rm -rf profile.0*
rm -rf tautrace.0*
make clean
make
ROCM_VERSION=`cat ${ROCM_PATH}/.info/version | head -1 | cut -f1 -d'-' `
result=`echo ${ROCM_VERSION} | awk '$1>6.1.9'` && echo $result
if [[ "${result}" ]]; then
   mpirun -n 2 tau exec -rocm -T rocm, rocprofsdk ./Jacobi hip -g 2 1
else
   mpirun -n 2 tau exec -T rocm,roctracer,rocprofiler ./Jacobi hip -g 2 1
fi
1s
pprof
```

Full code at: https://github.com/amd/HPCTrainingExamples/blob/main/tests/tau exec check.sh

Jacobi example (MPI+HIP) – Part 1/2

Using tau_exec with ROCm options (MPI is included unless the serial is declared)
 srun -n 2 --gpus 2 tau_exec -T rocm,rocprofiler,roctracer -rocm ./Jacobi_hip -g 2 1

Partial output of *pprof* command:

NODE 0	;CONTEXT 0;THREA	ND 0:				
%Time		Inclusive otal msec	#Call		Inclusive usec/call	
100.0			1	1	1942701	.TAU application
99.6	25	1,935	1	14020	1935195	.IAU application taupreload_main MPI_Init() hipMemcpy roctx: Jacobi_t::Jacobi_t::Top Level Init roctx: CreateMesh::Init hipMemset MPI Finalize() roctx: InitializeData::Init
32.5	630	630	1	1	630890	MPI_Init()
31.6	613	613	1005	Θ	611	hipMemcpy
18.8	0.059	365	1	3	365638	roctx: Jacobi_t::Jacobi_t::Top Level Init
15.6	0.138	303	1	6	303459	roctx: CreateMesh::Init
12.8	248	248	1	Θ	248057	hipMemset
7.6	147	147	1	2	147517	MPI_Finalize()
3.1	48	60	1	10	60800	roctx: InitializeData::Init
2.8	54	54	2	0	27380	hipStreamCreate hipStreamSynchronize roctx: Halo D2H::Halo Exchange
2.5	49	49	2000	0	25	hipStreamSynchronize
2.1	1	40	1000	2000	40	roctx: Halo D2H::Halo Exchange
1.7	2	33	1000	3000	34	roctx: MPI Exchange::Halo Exchange hipLaunchKernel MPI_Waitall()
1.5	29	29	5002	0	6	htpLaunchKernet
1.5	28	28	1000	9	29	MPI_Waitall()
1.2	1	22	1000	2000	23	roctx: Halo HZD::Halo Exchange
1.1	3	21	1003	3009	21	roctx: Halo H2D::Halo Exchange MPI_Allreduce() MPI Collective Sync
0.9	18	18	1005	9	18	MPI Collective Sync
0.6	10	10 7	2000	0	5	hipEventRecord hipMemcpyAsync
0.4	7	4	1000	0 0	/	n tpmemcpyAs ync
0.2	4	2		0	3	hipMemcpy2DAsync
0.1 0.1	2 2	2	1000 1001	9	3	MPI_Isend() hipDeviceSynchronize
0.1	1	1	7	9		hipMalloc
0.1	1	1		3		roctx: ApplyTopology::Init
0.1	0.22	1	1 1	3		MPI Comm dup()
0.1	1	1	1000	9		hipEventElapsedTime
0.1	0.983	0.983	2010	9	Ä	hipPointerGetAttributes
0.0	0.637	0.583	1000 2010 1000	9 9	1	MPI_Irecv()
0.0	0.525	0.525	3	9		hipHostMalloc
0.0	0.131			3	310	MPI_Allgather()
0.0	0.163	0.163	1 2	3 0	82	pthread join
0.0	0.012	0.012	2		6	hipGetDeviceCount
0.0	0.007	0.007				hinEventCreate
0.0	0.004	0.004	2 3	O	1	MPI_Barrier()
0.0		0.004	1	9	4	hipSetDevice
0.0	0.001	0.004 0.001	1	Ō	1	hipInit
			1	Θ		MDT Comm month()
Ofci.	21-23, 202	5 ₀	1	Θ		MPI_Comm_size() AMD @ Tsuki

- On the top of the output you see Node 0 (means first MPI process) and the thread 0 is always on the CPU, the names are about TAU naming and not the real processes
- Then you can see the percentage of the time, the exclusive duration and other information for all the calls from the code, MPI, and HIP API

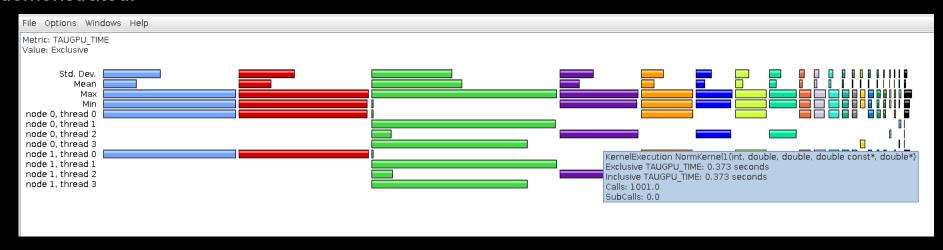
Jacobi example (MPI+HIP) – Part 2/2

- If you continue reading the output of the *pprof*, we can see the data transfers on thread 1 and 3, these are streams as the code uses stream
- Thread 2 is the actual process on the GPU, we can see the names of the kernels called, duration and other information

NODE 0;0	CONTEXT 0;THR	READ 1:				
%Time	Exclusive msec	Inclusive total msec	#Call	#Subrs	Inclusive usec/call	Name
100.0	877	890	1	1005		.TAU application
1.1 0.3	9 2	9 2	4 1001	0 0		CopyHostToDevice CopyDeviceToHost
0.3	2	2	1001	0	3	copyber (ceronos (
NODE 0;0	CONTEXT 0;THR	READ 2:				
%Time		Inclusive	#Call		Inclusive	Name
	msec	total msec			usec/call	
100.0	94	773	1	5002	773552	.TAU application
48.2	372	372	1001	0		KernelExecution No
21.8 16.5	168 127	168 127	1000 1000	9 9		KernelExecution Ja KernelExecution Lo
1.0	8	8	1000		8	KernelExecution Ha
0.3	2	2	1001	Θ		KernelExecution No
NODE 0;0	CONTEXT 0;THR	READ 3:				
%Time	Exclusive	Inclusive	#Call	#Subrs	Inclusive	Name
	msec	total msec			usec/call	
100.0	744	771	1	2000	771690	.TAU application
3.1	23	23	1000	0		CopyDeviceToDevice
0.5	3	3	1000	0	4	CopyHostToDevice

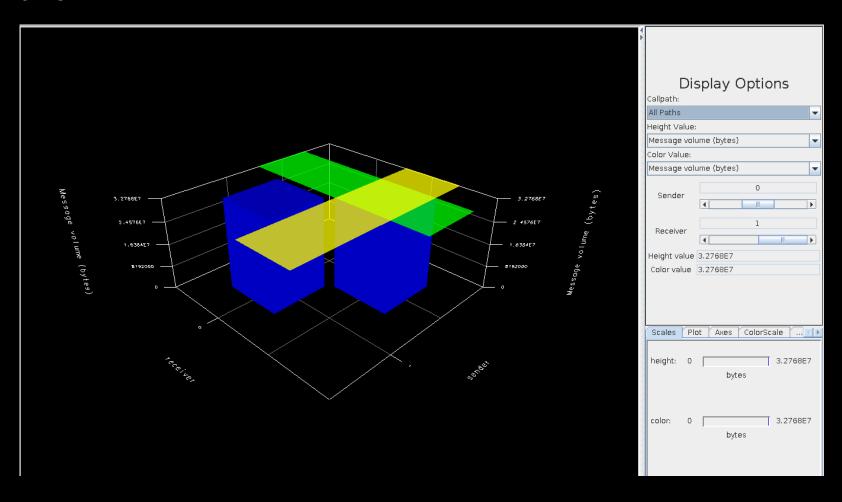
Jacobi visualization with Paraprof

- Pack the profiling data to a file:
 paraprof --pack jacobi.ppk
- Then execute the paraprof tool paraprof jacobi.ppk
- Each color represents a different routine/API call, each row maps to the information on the left, mean/min/max/stdev values across all the threads, when you go the mouse over a color, it pop ups the name of the event as it is demonstrated.



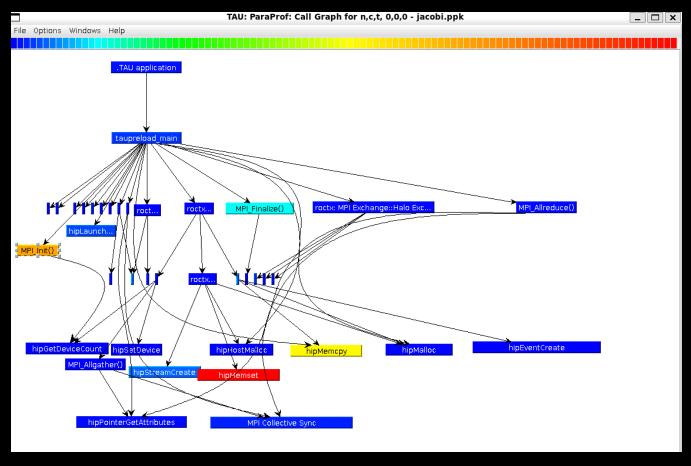
Jacobi - Communication Matrix

Windows -> 3D Communication Matrix



Jacobi - Callgraph

- Windows -> Thread -> Callgraph and select a thread, we chose the CPU thread (0)
 - The color indicates which calls take more time than others (blue color) as also you can see the call dependencies



Jacobi - Tracing

- Activate tracing: export TAU_TRACE=1
- Now we have one TRC file per thread and one EDF file per MPI process
- Merge all the files with the command: tau_treemerge.pl now there are two files named tau.trc and tau.edf

• Create a json file to visualize with perfetto: tau_trace2json tau.trc tau.edf -chrome -ignoreatomic -o jacobi.json



Porting an MPI CPU application

Identify hot loops for CPU code using TAU with PDT

Name	Exclusive TIME ▽	Inclusive TIME	Calls	Child Calls
Loop: SETUPMOD::FELLI_MESHREF [{setupmod.f90} {2392,5}-{2534,10}]	6.072	8.979	896	395,065 📤
- MATHMOD::HYBRD [{mathmod.f90} {2707,12}] [THROTTLED]	2.045	2.547	254,745	1,366,010
- Loop: MOMMOD::VIS3D [{mommod.f90} {1842,1}-{2654,16}]	0.983	1.02	10	1,330
- Loop: MGCCMOD::SMOOTH [{mgccmod.f90} {2923,7}-{2943,12}]	0.561	0.561	5,771	0 =
MATHMOD::HYBRD1 [{mathmod.f90} {3255,12}] [THROTTLED]	0.349	2.897	289,683	254,745
MGCCMOD::RESTRICT [{mgccmod.f90} {4706,12}]	0.27	0.393	4,179	16,716
Loop: GEOMETRYMOD::GRDF3D [{geometrymod.f90} {966,5}-{1291,10}]	0.254	0.384	54	676,999
Loop: GODSPLIT3D::Z VEL CALC CONS [{godmod.f90} {1625,1}-{1653,6}]	0.232	0.241	10	100,001
Loop: GODSPLIT3D::YVEL_CALC_CONS [{godmod.f90} {1162,1}-{1189,6}]	0.222	0.222	10	О
Loop: GODSPLIT3D::X VEL CALC CONS [{godmod.f90} {712,1}-{740,6}]	0.222	0.222	10	o
Loop: MGCCMOD::SMOOTH [{mgccmod.f90} {2395,7}-{2429,12}]	0.18	0.18	199	0
Loop: MGCCMOD::PROLONGATE [{mgccmod.f90} {6393,3}-{6408,8}]	0.131	0.131	4,179	0
Loop: MGCCMOD::RESTRICT [{mgccmod.f90} {5380,9}-{5421,14}]	0.122	0.122	4,179	0
Loop: GODMOD::GODSPLIT3D [{godmod.f90} {349,1}-{632,13}]	0.102	1.36	10	120
MPI File write at all()	0.1	0.1	48	145
Loop: GODSPLIT3D::Z VEL_CALC_CONS [{godmod.f90} {1678,1}-{1693,6}]	0.077	0.077	10	0
Loop: GODSPLIT3D::Z VEL CALC_CONS [{godmod.f90} {1658,1}-{1673,6}]	0.076	0.076	10	ő
Loop: GODSPLIT3D::Y VEL CALC CONS [{godmod.f90} {1213,1}-{1227,6}]	0.075	0.075	10	0
Loop: GODSPLIT3D::X VEL CALC CONS [{godmod.f90} {764,1}-{778,6}]	0.075	0.075	10	n i
Loop: GODSPLIT3D::Y VEL CALC CONS [{qodmod.f90} {1194,1}-{1208,6}]	0.075	0.075	10	ő
Loop: GODSPLIT3D::X VEL CALC CONS [{godmod.f90} {745,1}-{759,6}]	0.075	0.075	10	o o
Loop: SETUPMOD::SETUP3D [{setupmod.f90} {407.3}-{865.8}]	0.074	9.116	10	109,377
Loop: SETUPMOD::SETUP3D [{setupmod.fs0} {914,5}-{922,10}]	0.074	0.117	1	144,209
MATHMOD::R1UPDT [{mathmod.f90} {3797,12}] [THROTTLED]	0.072	0.105	100,001	400,004
BCMOD::BC VECTOR [{bcmod:f90} {1813,12}]	0.054	0.065	1,294	9,078
GEOMETRYMOD::GRDF3D [{geometrymod.f90} {155,12}]	0.034	0.437	54	486
MATHMOD::QFORM [{mathmod.f90} {3391,12}] [THROTTLED]	0.045	0.437	100,001	300,003
FS3DMOD::FS3D INIT [{fs3d.f90} {1684,12}]	0.043	9.519	100,001	40
	0.043	0.042	21	40
Loop: MGCCMOD::DISCRETIZE [{mgccmod.f90} {7158,3}-{7178,8}]	0.042	0.042	138	0
MPLFile_write_at()	0.037	4.547	138	170
FS3DMOD::EULER_EXPL [{fs3d.f90} {8154,12}]	0.034			200,002
MATHMOD::QRFAC [{mathmod:f90} {3492,12}] [THROTTLED]		0.083	100,001	
MATHMOD::DOGLEG [{mathmod.f90} {2267,12}] [THROTTLED]	0.034	0.051	100,001	164,917
MATHMOD::R1MPYQ [{mathmod.f90} {3686,12}] [THROTTLED]	0.033 0.033	0.053 0.056	100,001	200,002
GEOMETRYMOD::INTERFACE_HEIGHT_TOP [{geometrymod.f90} {4730,12}] [THROTTLED]				200,002
SETUPMOD::FELLI_MESHREF [{setupmod.f90} {2252,12}] [THROTTLED]	0.029	0.038	100,001	100,001
Loop: MATHMOD::FDJAC1 [{mathmod.f90} {2639,6}-{2658,11}] [THROTTLED]	0.028	0.032	100,001	41,460
Loop: SURFACETENSIONMOD::STEN3D [{surfacetensionmod.f90} {581,1}-{665,6}]	0.027	0.027	11	0
Loop: MATHMOD::QRFAC [{mathmod.f90} {3611,3}-{3682,8}] [THROTTLED]	0.025	0.026	100,001	11,124
.TAU application	0.024	14.322	1	1
Loop: MGCCMOD::SMOOTH [{mgccmod.f90} {2904,7}-{2919,12}]	0.024	0.024	5,771	0
Loop: MGCCMOD::SMOOTH [{mgccmod.f90} {3011,7}-{3027,12}]	0.023	0.023	5,771	0
SETUPMOD::VORBELEGUNGSFUNKTIONEN [{setupmod.f90} {5382,12}] [THROTTLED]	0.022	0.061	100,001	100,001
MATHMOD::FDJAC1 [{mathmod.f90} {2528,12}] [THROTTLED]	0.022	0.054	100,001	100,001
MATHMOD::RGAUSS [{mathmod.f90} {106,30}] [THROTTLED]	0.021	0.042	100,001	100,001
Loop: MATHMOD::QRFAC [{mathmod.f90} {3594,3}-{3596,8}] [THROTTLED]	0.021	0.022	100,001	11,124
Loop: MGCCMOD::RAP [{mgccmod.f90} {6609,3}-{6650,8}]	0.021	0.021	126	0
MGCCMOD::BC_ALL [{mgccmod.f90} {4443,12}]	0.021	0.021	199	199
MPI_File_open()	0.02	0.02	10	0 -
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Fortran+OMP offloading (Nekbone)

- Execute the application and declare to tau_exec to profile ROCm events (MPI is activated by default, otherwise define SERIAL)
- srun –n1 tau_exec -T rocm,rocprofiler,roctracer -rocm ./nekbone data

File Option	ns Windows	Help					
Metric: TAL Sorted By: Units: seco	Exclusive						
%Tota	l Time	Exclusive	Inclusive	#Calls	#Child Calls	Inclusive/Call	Name
	100.0	0.687	0.993	1	2614	0.993	.TAU application
	7.7	0.076	0.076	200	0	3.8E-4	omp offloading 25c66b0a c60lfl35 ax acc l46l.kd
	7.2	0.072	0.072	602	0	1.2E-4	omp_offloading_25c66b0a_c601flc1_glsc3_accl1447.kd
	7.1	0.071	0.071	200	0	3.5E-4	omp_offloading_25c66b0a_c601f135_ax_accl496.kd
	3.8	0.038	0.038	600	0	6.3E-5	_omp_offloading_25c66b0a_c601flc1_add2s2_accl1483.kd
	1.3	0.012	0.012	200	0	6.2E-5	_omp_offloading_25c66b0a_c601flc1_add2s1_accl1467.kd
	1.0	0.01	0.01	200	0	4.9E-5	_omp_offloading_25c66b0a_c601flcd_dssum_accl458.kd
	0.9	0.009	0.009	200	0	4.7E-5	_omp_offloading_25c66b0a_c601flc1_copy_accl1425.kd
	0.9	0.009	0.009	200	0	4.5E-5	omp_offloading_25c66b0a_c601flcd_dssum_accl486.kd
	0.8	0.008	0.008	200	0	3.8E-5	omp_offloading_25c66b0a_c601f135_maskit_accl655.kd
	0.0	1.6E-4	1.6E-4	4	0	3.9E-5	omp_offloading_25c66b0a_c601f17d_gather_double_add_l148.kd
	0.0	1.3E-4	1.3E-4	4	0	3.1E-5	omp_offloading_25c66b0a_c60lfl7d_scatter_double_l148.kd
	0.0	1.0E-4	1.0E-4	2	0	5.0E-5	omp_offloading_25c66b0a_c60lflcl_rzero_accl1393.kd
	0.0	6.0E-6	6.0E-6	2	0	3.0E-6	omp_offloading_25c66b0a_c601f17d_init_double_l148.kd

HPCToolkit



HPCToolkit: Overview

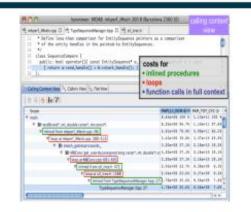
- Not an AMD profiler developed for 20+ years, mainly at Rice University
- HPCToolkit suite of tools for tracing, profiling and analyzing parallel programs
- Combine sampling data with a static analysis of the program structure for loops and inline functions
- Present top-down, bottom-up and flat views of calling context tree and time-sequence trace view
- Low overhead, easy to use, interactive analysis with hpcviewer
- No instrumentation or recompilation needed, as long as "-g" compiler flag is used
- Supports threads (pthreads, OpenMP®), MPI, hybrid (MPI+threads), and GPUs (AMD, Intel®, NVIDIA)

Installing HPCToolkit

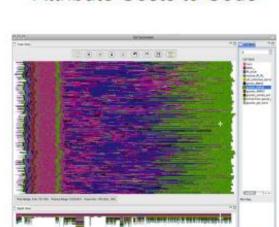
- HPCToolkit integrated suite of tools for measurement and analysis of program performance on computers ranging from multicore desktop systems to GPU-accelerated supercomputers (Rice University)
- Installation from <u>Gitlab</u> repo using **Meson**:

```
pipx install 'meson>=1.3.2'
export PATH=$HOME/.local/bin:$PATH
git clone https://gitlab.com/hpctoolkit/hpctoolkit.git
git checkout 0bfc9cd58da73e3e5b973d754677ebfe2b0da55d
cd hpctoolkit
export CMAKE PREFIX PATH=$ROCM PATH:$CMAKE PREFIX PATH
meson setup -Drocm=enabled --prefix=${HPCTOOLKIT PATH}
            --libdir=${HPCTOOLKIT PATH}/lib build
  build
cd
meson compile
meson install
```

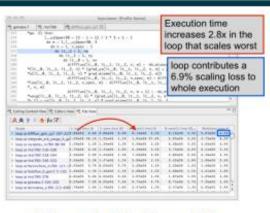
HPCToolkit Capabilities at a Glance



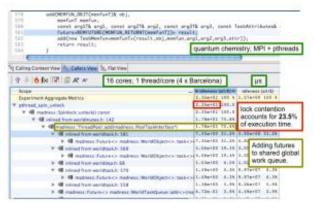
Attribute Costs to Code



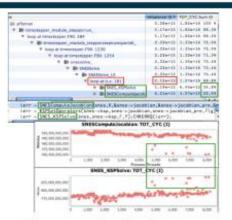
Analyze Behavior over Time



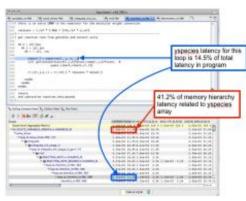
Pinpoint & Quantify Scaling Bottlenecks



Shift Blame from Symptoms to Causes



Assess Imbalance and Variability

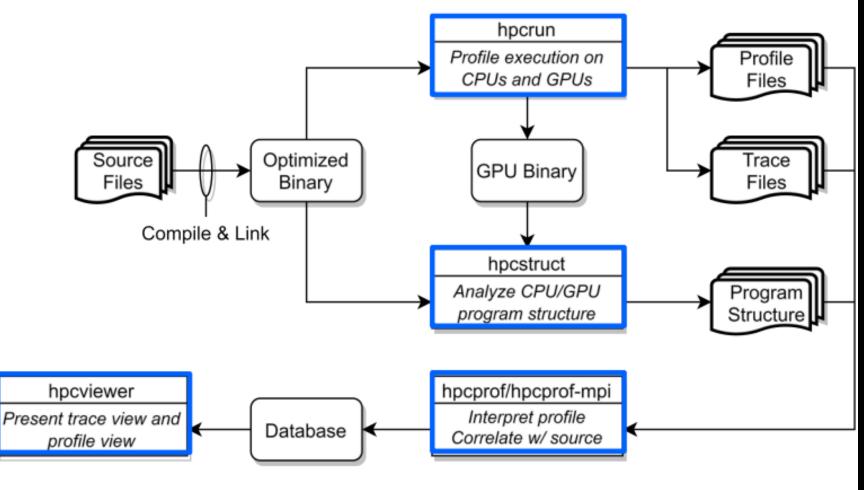


Associate Costs with Data





HPCToolkit's Workflow for GPU-accelerated Applications





HPCToolkit examples in HPCTrainingExample repo

hpcrun

```
# This test checks that hpcrun
# runs without errors

module load rocm

module load hpctoolkit
REPO_DIR="$(dirname "$(dirname "$(readlink -fm "$0")")")"
pushd ${REPO_DIR}/HIP/Stream_Overlap/0-Orig/
rm -rf build_for_test
mkdir build_for_test; cd build_for_test
cmake ../
make -j

hpcrun -e CPUTIME -e gpu=amd -t ./compute_comm_overlap 2
ls hpctoolkit-compute_comm_overlap-measurements*
```

hpcstruct

```
# This test checks that hpcstruct
# runs without errors

module load rocm

module load hpctoolkit
REPO_DIR="$(dirname "$(readlink -fm "$0")")")"
pushd ${REPO_DIR}/HIP/Stream_Overlap/0-Orig/
rm -rf build_for_test
mkdir build_for_test; cd build_for_test
cmake ../
make -j

hpcrun -e CPUTIME -e gpu=amd -t ./compute_comm_overlap 2
hpcstruct hpctoolkit-compute_comm_overlap-measurements*
ls hpctoolkit-compute_comm_overlap-measurements*
```

hpcprof

```
# This test checks that hpcprof
# runs without errors

module load rocm

module load hpctoolkit
REPO_DIR="$(dirname "$(dirname "$(readlink -fm "$0")")")"
pushd ${REPO_DIR}/HIP/Stream_Overlap/0-Orig/
rm -rf build_for_test
mkdir build_for_test; cd build_for_test
cmake ../
make -j

hpcrun -e CPUTIME -e gpu=amd -t ./compute_comm_overlap 2
hpcstruct hpctoolkit-compute_comm_overlap-measurements*
hpcprof hpctoolkit-compute_comm_overlap-measurements*
ls hpctoolkit-compute_comm_overlap-database*
```

Full code at: https://github.com/amd/HPCTrainingExamples/blob/main/tests/

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