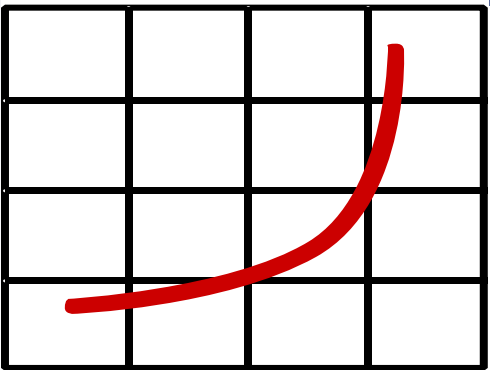


How to run SPEC benchmarks?

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spec

- Hands-On: Cluster login
- Overview of System Requirements
- Configuration Files – The bad, the good & the ugly
 - SPEC ACCEL: Investigating OpenACC Performance
 - runspec & run rules
 - From base to peak runs
 - Different compilers
 - SPEC ACCEL: Adaptions for OpenCL Benchmarks
 - SPEC OMP2012: Host Benchmarking
- Hands-On: Run your benchmark

Cluster login

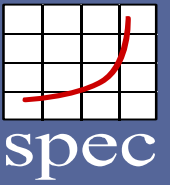
- See instructions on handouts
- Follow interactive demo

Interactive demo time!

- We present SPEC ACCEL OpenACC config files
- More slides for OpenCL and OpenMP attached
- Opportunity to follow instructions interactively (also see handouts)
- Later: run benchmarks
 - Choose from OpenACC, OpenCL or OpenMP (whatever you like)
 - MPI runs take very long – get prepared results
- Advertisement: interpret results after break

- Hands-On: Cluster login
- **Overview of System Requirements**
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System Requirements



- Different benchmarks suites – different requirements
 - SPEC ACCEL¹, SPEC OMP2012², SPEC MPI2007³
- Supported operating systems: AIX, Linux, MacOS, Solaris, Windows (except very old Windows)
 - Please do not use Windows/Unix compatibility products
- Compatible processors
 - CPU
 - GPU
 - APU
 - Xeon Phi

¹ <https://www.spec.org/accel/docs/system-requirements.html>

² <https://www.spec.org/omp2012/Docs/system-requirements.html>

³ <https://www.spec.org/mpi2007/Docs/system-requirements.html>

- Memory requirements
 - OpenMP: 28GB for the whole system
 - MPI: 1GB/rank (medium size) and 2GB (large size)
 - ACCEL: 4GB of host + 2GB of device
 - Otherwise, you are measuring your paging file, not your system
- Disk space requirements
 - OpenMP: 8GB
 - MPI: 10GB (medium), 17GB (large, big endian), 24GB (large, little endian)
 - ACCEL: 9GB
- Support of compilers
 - C99, C++98 and Fortran-95 compilers + MPI library for SPEC MPI 2007

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Tools provided to ensure consistent operation of benchmarks across variety of platforms¹

- **specmake**
 - GNU make to build benchmarks
- **runspec**²
 - Primary tool in the suite
 - Used to build the benchmarks, run them, and report their results
 - Config file needed for usage (with detailed instructions about how to build and run the benchmarks)
- And more

¹<https://www.spec.org/accel/docs/tools-build.html>

²<https://www.spec.org/accel/docs/runspec.html>

runspec - Run SPEC benchmarks

```
$> cd $SPEC
$> . ./shrc
$>
$> runspec --config=tutSC15-openacc-pgi --tune=base,peak 350
```

- Base & peak runs

Use config file
myopenaccconf

- Run selected benchmark: 350
- Or: whole suite, e.g. openacc

- Compile benchmarks
- Run benchmarks
- Obey to **SPEC run rules**¹
 - Tester supplies compiler & system
 - No code modifications of provided sources allowed
 - Restrictions for base & peak runs
 - Data set sizes pre-specified (test, train, ref)²

¹ <https://www.spec.org/accel/docs/runrules.html>,
<https://www.spec.org/omp2012/docs/runrules.html>,
<https://www.spec.org/mpi/docs/runrules.html>

² For MPI2007: medium and large data sets

- Contain instructions for
 - building benchmarks
 - running them
 - description of system under test

Key for reproducibility!

Use Case:

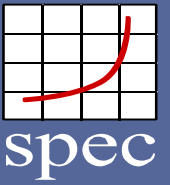
How to investigate OpenACC performance? The bad, the good & the ugly

- How to write a config file? – The bad
 - Often start off using a config file that someone else has previously written¹
 - E.g. directory `$SPEC/config/` or ACCEL result submissions similar to your system
 - Write your own²

¹ https://www.spec.org/accel/docs/runspec.html#about_config

² <https://www.spec.org/accel/docs/config.html>

Structure of Config Files



Header section

- 1st section prior to names section
- Usually runspec flags

```
iterations      = 1
# Tuning levels: base, peak
tune            = base
# Dataset size: test, train, ref
size           = ref
# Environment variable will be set using "ENV_*"
env_vars       = 1
# Output format: all, pdf, text, html and so on
output_format  = text
flagsurl       = ${top}/result/pgi_flags.xml
teeout         = yes

# Run benchmarks according to your specific system config
# The variable "command" is the command used by spec
submit = aprun -n 1 -N 1 -q $command
```

```
#####
# Compiler information
#####
```

Named section

- Begins w/ “section marker”
- One- to four-part string of the form
benchmark[,...]=tuning[,...]
=extension[,...]=machine[,...]:

```
openacc=base=default=default:
OPTIMIZE      = -fast -Mfpelaxed
FOPTIMIZE     = -acc -ta=tesla:cc35,cuda6.5
COPTIMIZE     = -acc -ta=tesla:cc35,cuda6.5
```

```
#####
# Portability flags for each benchmark
# Following flag should not have any impact on performance.
#####
116.histo=default=default=default:
PORTABILITY   = -DSPEC_LOCAL_MEMORY_HEADROOM
```

```
118.cutcp=default=default=default:
CPORTABILITY += -D__GNUC__

359.miniGhost=default=default=default:
EXTRA_LDFLAGS += -Mnomain
```

```
#####
# Peak
#####
350.md=peak=default=default:
FOPTIMIZE     = -acc -ta=tesla:cc35,cuda5.5,maxregcount:54
```

```
363.swim=peak=default=default:
FOPTIMIZE     = -acc -ta=tesla:cc35,cuda5.5,pin
```

```
#####
# Hardware and software information for the machine under test.
# This information will be extracted for a reportable run.
# An example configuration can be copied from the website
# https://www.spec.org/accel/results/accel_acc.html
#####
```

```
company_name   = SPEC Tutorial Company
test_sponsor   = SPEC Tutorial Sponsor
tester        = SPEC Tutorial Tester
license_num    = SPEC Tutorial License
machine_name   = SPEC Tutorial Machine
hw_vendor      = Cray
hw_avail       = Apr-2013
hw_cpu         = AMD Opteron Processor 6276
hw_cpu_mhz     = 2300
hw_cpu_max_mhz = 3200
hw_cpu_char000 = AMD Turbo CORE Technology up to 3.2GHz,
hw_cpu_char001 = Turbo CORE off
```

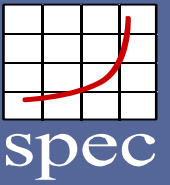
```
hw_disk        = None
hw_fpu         = Integrated
hw_memory      = 32 GB (4 x 8 GB 2Rx4 PC3L-12800R-11, ECC)
hw_model       = Cray XK7
hw_ncores      = 16
hw_nchips      = 2
hw_ncoresperchip = 16
hw_nthreadspercore = 1
hw_ncpuorder   = 1 chip
hw_other       = None
hw_ocache      = None
hw_pcache      = 32 KB I + 16 KB D on chip per core
hw_scache      = 16 MB I+D on chip per chip, 2 MB shared / 2 cores
hw_tcache      = 16 MB I+D on chip per chip, 8 MB shared / 8 cores
hw_avail       = Feb-2015
sw_compiler000 = PGI Accelerator Fortran/C/C++ Server,
sw_compiler001 = Release 15.3
sw_file        = NFSv3 (IBM N5500 NAS) over Gb ethernet
sw_os000       = SUSE Linux Enterprise Server 11 (x86_64),
sw_os001       = Cray Linux Environment 4.2, Kernel
hw_os002       = 2.6.32.59-0.7.1_1.0402.7496-cray_gem_c
sw_other       = NVIDIA CUDA 5.5.20
sw_state       = Multi-user, run level 3
sw_base_ptrsize = 64-bit
hw_accel_connect = PCIe 2.0 16x
hw_accel_desc000 = NVIDIA Tesla K20m GPU, 2496 CUDA cores,
hw_accel_desc001 = 706MHz, 5 GB GDDR5 RAM
hw_accel_ecc     = yes
hw_accel_model   = Tesla K20
```

MD5 section

- Automatically-generated

```
#####
# MD5 section. It will be created by SPEC automatically.
# It is used by SPEC to check whether an executable if
# available is created using the current compiler and flags
# settings.
#####
```

SPEC ACCEL Config File (1/11)



config file: \$SPEC/config/tutSC15-openacc-pgi.cfg

```
#####
```

```
# The header section of the config file. Must appear  
# before any instances of "default="
```

```
#####
```

```
# what to do: build, validate = build + run  
action      = validate
```

```
# Number of iterations of a test  
iterations   = 1
```

```
# Tuning levels: base, peak  
tune         = base
```

```
# Dataset size: test, train, ref  
size         = ref
```

- **build**: compile benchmarks
- **validate**: benchmarks are built if necessary, run and reports are generated

- How many times to run each benchmarks
- e.g. for reportable run = 3

- **base**: flags for all benchmarks the same
- **peak**: set of optimizations individually selected for that benchmark

- Data set sizes (from small to big): test, train, ref
- e.g. test for debugging new set of compilation options

SPEC ACCEL Config File (2/11)

config file: \$SPEC/config/tutSC15-openacc-pgi.cfg

```
#####  
# The header section of the config file. Must appear  
# before any instances of "default="  
#####
```

- Environment settings
- ENV_VAR = ...
- Apply to build phase → rebuild if any changes

```
# Environment variable will be set using "ENV_*", see the next section
```

```
env_vars      = 1
```

- Different output formats possible
- In \$SPEC/results

```
# Output format: all, pdf, text, html and so on
```

```
output_format = text
```

```
flagsurl      = ${top}/result/pgi_flags.xml
```

```
teeout        = yes
```

- Description of portability & tuning options ("Flags File")
- Information on syntax of flags and their meanings
- Needed for valid reports

Displays the build commands to screen

SPEC ACCEL Config File (3/11)

```
# How to run the benchmarks according to your specific system configuration
# The variable "command" is the command used by spec
submit = aprun -n 1 $command
```

- **How to execute the benchmarks**
- Use `$command` for SPEC command
- Preferred to assign work to processors
 - May place benchmarks on desired processors or benchmark memory on a desired memory unit
 - Especially needed for MPI runs
 - Here: run job on one node
- Can be used to change the run time environment

```
submit = export ENV_VAR=...; ...
```

→ no rebuild if any changes occur

SPEC ACCEL Config File (4/11)

```
#####
# Software information
#####
```

```
# Compilers. Using PGI compiler for example
default=default=default=default:
CC          = pgcc
CXX         = pgc++
FC          = pgfortran
```

```
#####
# Base
#####
```

```
openacc=base=default=default:
OPTIMIZE   = -fast -Mfprelaxed
FOPTIMIZE  = -acc -ta=tesla:cc35,cuda5.5
COPTIMIZE  = -acc -ta=tesla:cc35,cuda5.5
```

Named section

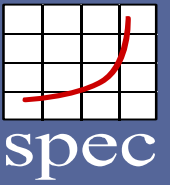
```
benchmark[,...]=tuning[,...]
=extension[,...]=machine[,...]:
```

Base

- Common set of optimizations & environment settings for all benchmarks
- “baseline”
 - single set of switches
 - single-pass make process
 - high degree of portability, safety, performance
- Must adhere to strict rules
 - e.g. same compiler for all modules of a given language
 - All flags, options must be the same
 - More rules (base & peak) on names, library substitutions, data type sizes, source code changes

- Define PGI compiler for OpenACC compilation
- Define PGI GPU flags

SPEC ACCEL Config File (5/11)



```
#####  
# Portability flags for each benchmark  
# Following flag should not have any impact on performance.  
#####
```

```
116.histo=default=default=default:  
PORTABILITY = -DSPEC_LOCAL_MEMORY_HEADROOM
```

```
118.cutcp=default=default=default:  
CPORTABILITY += -D__GNUC__
```

```
359.miniGhost=default=default=default:  
EXTRA_LDFLAGS += -Mnomain
```

Portability flags

- Allowed if benchmark cannot be built and execute correctly w/o these flags
- Must be performance neutral
- Base rules except portability flags, i.e. flags may differ from one benchmark to another (even in base)
- Requirements
 - Provided over PORTABILITY flag
 - Must be approved by SPEC HPG committee

SPEC ACCEL Config File (6/11)

```
#####
# Hardware and software information for the machine under test.
# This information will be extracted for a reportable run.
# An example configuration can be copied from the website
# https://www.spec.org/accel/results/accel_acc.html
#####
```

```
company_name      = SPEC Tutorial Company
test_sponsor      = SPEC Tutorial Sponsor
tester            = SPEC Tutorial Tester
license_num       = SPEC Tutorial License
machine_name      = SPEC Tutorial Machine
```

```
hw_vendor         = Cray
hw_avail          = Apr-2013
hw_cpu            = AMD Opteron Processor 6276
hw_cpu_mhz        = 2300
hw_cpu_max_mhz    = 3200
```

HW & SW description

- Needed only for reportable runs
- runspec tools captures information in submission file
- Very detailed information
- More: next tutorial section

Information on host configuration, e.g. CPU

SPEC ACCEL Config File (7/11)

```
hw_cpu_char      = AMD Turbo CORE Technology up to 3.2GHz, Turbo CORE off
hw_disk          = None
hw_fpu           = Integrated
hw_memory        = 32 GB (4 x 8 GB 2Rx4 PC3L-12800R-11, ECC)
hw_model         = Cray XK7
hw_ncores        = 16
hw_nchips        = 2
hw_ncoresperchip = 16
hw_nthreadspercore = 1
hw_ncpuorder     = 1 chip
hw_other         = None
hw_ocache        = None
hw_pcache        = 32 KB I + 16 KB D on chip per core
hw_scache        = 16 MB I+D on chip per chip, 2 MB shared / 2 cores
hw_tcache        = 16 MB I+D on chip per chip, 8 MB shared / 8 cores
```

SPEC ACCEL Config File (8/11)

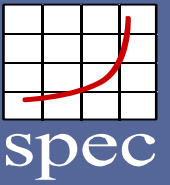
```
sw_avail      = Feb-2015
sw_compiler   = PGI Accelerator Fortran/C/C++ Server, Release 15.3
sw_file       = NFSv3 (IBM N5500 NAS) over Gb ethernet
sw_os000      = SUSE Linux Enterprise Server 11 (x86_64), Cray Linux Environment 4.2
sw_os001      = Kernel 2.6.32.59-0.7.1_1.0402.7496-cray_gem_c
sw_other      = NVIDIA CUDA 5.5.20
sw_state      = Multi-user, run level 3
sw_base_ptrsize = 64-bit
```

Information on software
configuration, e.g. compilers

```
hw_accel_connect = PCIe 2.0 16x
hw_accel_desc    = NVIDIA Tesla K20m GPU, 2496 CUDA cores, 706MHz, 5 GB GDDR5 RAM
hw_accel_ecc     = yes
hw_accel_model   = Tesla K20
hw_accel_name    = NVIDIA Tesla K20
hw_accel_type    = GPU
hw_accel_vendor  = NVIDIA
sw_accel_driver  = NVIDIA UNIX x86_64 Kernel Module 319.82
```

Information on accelerator
configuration, e.g. device
vendor

SPEC ACCEL Config File (9/11)



```
#####  
# MD5 section. It will be created by SPEC automatically.  
# It is used by SPEC to check whether an executable if available is created using  
# the current compiler and flags settings.  
#####
```

MD5 section

- Automatically generated by SPEC tools
- Used to check whether an executable is created using the current settings

SPEC ACCEL Config File (10/11)

- Modifying the config file – The good
 - Once you have a config file that runs on your system, it is easy to modify it
 - E.g. `peak` optimizations for better performance

```
#####
# Peak
#####

350.md=peak=default=default:
FOPTIMIZE      = -acc -ta=tesla:cc35,cuda5.5,maxregcount:54

363.swim=peak=default=default:
FOPTIMIZE      = -acc -ta=tesla:cc35,cuda5.5,pin
```

Peak

- Set of optimizations individually selected for each benchmark
 - e.g. different compilers, flags
- Called “aggressive compilation”

- `maxregcount`: sets register spilling to 54
- `pin`: pinned memory usage

- `md` benchmark: computational-intensive
- `swim` benchmark: many data transfers between CPU & GPU

SPEC ACCEL Config File (11/11)

- Rewriting the config file – The ugly
 - Major changes in a config file (e.g. compiler) need more adaption

```
# Compilers. Using Cray compiler for example
default=default=default=default:
CC          = cc
CXX         = CC
FC          = ftn
#####
# Base
#####
openacc=base=default=default:
OPTIMIZE    = -O2
FOPTIMIZE   = -h acc,noomp -em -fpic -dynamic
COPTIMIZE   = -h pragma=acc -h nopragma=omp -fpic -dynamic
#####
# new portability flags; new peak flags
#####
```

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SPEC ACCEL OpenCL – Device Selection

- SPEC ACCEL OpenCL benchmarks are compatible with multiple devices
- These devices include: CPU, GPU, Xeon Phi and APU
- When multiple devices co-exist, SPEC ACCEL OpenCL allows user to specify the device for running benchmarks through the runspec option
`--device <device_number>`
- *device_number* can be a
 - number
 - name among CPU, GPU and ACCELERATOR

Example:

```
$> runspec --size train --platform NVIDIA --device 0 --config tutSC15-ocl-gnu.cfg lavamd
```


SPEC ACCEL OpenCL – Workgroup Sizes

- Some (not all) OpenCL benchmarks have the ability to change the workgroup size
- Changing workgroup size may improve the performance
- To accomplish this, SPEC provides define variables to change workgroup sizes
- Some of the benchmarks have multiple variables for different OpenCL kernels
- Only allowed in a peak run.

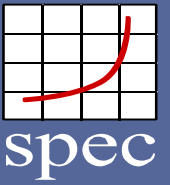
- For one benchmark:

- SPEC_ACCEL_WG_SIZE
 - sets all of the workgroup sizes to this value
- SPEC_ACCEL_WG_SIZE_n and SPEC_ACCEL_WG_SIZE_n_n
 - set specific values for a specific kernel and a specific dimension of a kernel, respectively

101.tpacf	0	0_0
103.stencil	0, 1	0_0, 0_1, 0_2
104.lbm	0, 1	0_0, 0_1, 0_2
112.spmv	0	0_0
114.mriq	0, 1	0_0, 1_0
116.Histo	0, 1, 2, 3	0_0, 1_0, 2_0, 3_0
117.bfs	0, 1	0_0, 1_0
120.kmeans	0, 1	0_0, 1_0
121.lavamd	0	0_0
122.cfd	0, 1, 2, 3, 4	0_0, 1_0, 2_0, 3_0, 4_0
123.nw	0	0_0
124.hotspot	0	0_0
125.lud	0	0_0
127.srad	0	0_0
128.heartwall	0	0_0

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SPEC OMP2012 Config File (1/4)



config file: \$SPEC/config/myopenmpconf.cfg

```
#####  
# Header section  
#####  
  
# what to do build  
action          =      validate  
# Number of iterations of a test  
iterations      =      1  
# Tuning levels: base, peak  
tune            =      base,peak  
# Dataset size: test, train, ref  
size            =      train  
# Number of used threads  
threads         =      16  
# Environment variable will be set using "env_vars", see the next section  
env_vars        =      1  
# Output format: all, pdf, text, html and so on  
output_format   =      text
```

OMP needs # threads

OR

runspec --config=myopenmpconf --threads=16 omp

Environment variables
important for NUMA
settings

SPEC OMP2012 Config File (2/4)

```
#####
# Compiler information
#####
# Compilers. Using Intel compiler for example
default=default=default=default:
CC  = icc
CXX = icpc
FC  = ifort
F77 = ifort
#####
# Portability flags for each benchmark
#####
350.md=default=default=default:
FPORTABILITY  = -free
367.imagick=default=default=default:
CPORTABILITY  = -std=c99
357.bt331=default=default=default:
PORTABILITY   = -mcmodel=medium
363.swim=default=default=default:
PORTABILITY   = -mcmodel=medium
```

Choose your OMP compiler

As always: portability flags should not impact performance

SPEC OMP2012 Config File (3/4)

```
#####
# Base
#####
default=base=default=default:
# Basic optimization flags for all benchmarksE
OPTIMIZE=-O3 -openmp -ipo -xCORE-AVX-I -no-prec-div
# Optimization flags for benchmarks written in C
COPTIMIZE=-ansi-alias
# Optimization flags for benchmarks written in C++
CXXOPTIMIZE=-ansi-alias
# Optimization flags for benchmarks written in Fortran
FOPTIMIZE=-align
# Environment variables at runtime
ENV_KMP_AFFINITY = compact
ENV_KMP_SCHEDULE = static,balanced
ENV_KMP_BLOCKTIME = infinite
ENV_KMP_LIBRARY = throughput
ENV_KMP_STACKSIZE = 500M
ENV_OMP_NESTED = FALSE
ENV_OMP_DYNAMIC = FALSE
```

Basic optimization flags

Environment variables,
e.g. on NUMA settings
or loop schedules

SPEC OMP2012 Config File (4/4)

```
#####
```

```
# Peak
```

```
#####
```

```
350.md=peak=default=default:
```

```
OPTIMIZE=-O3 -openmp -ipo -xCORE-AVX-I -ansi-alias -opt-malloc-options=1
```

```
FOPTIMIZE=-fp-model fast=2 -no-prec-div -no-prec-sqrt -align array64byte
```

Peak runs include more aggressive optimizations

```
363.swim=peak=default=default:
```

```
OPTIMIZE=-O3 -openmp -ipo -xCORE-AVX-I -no-prec-div -ansi-alias -opt-streaming-stores always  
-opt-malloc-options=4
```

```
ENV_KMP_AFFINITY=compact,0
```

```
#####
```

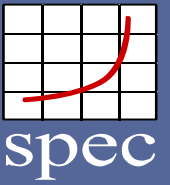
```
# Hardware information
```

```
#####
```

```
# MD5 section
```

```
#####
```

Summary - SPEC ACCEL & SPEC OMP2012



- SPEC ACCEL for accelerators
 - Including benchmarks implemented using OpenACC and OpenCL
 - Under development: benchmarks using OpenMP target
 - Don't need to specify #threads
 - Can specify an accelerator, if there are multiple
- SPEC OMP2012 for Host CPU
 - Benchmarks implemented using OpenMP for host
 - Need to specify #threads
 - Processor/core can be chosen using environment variable, such as ENV_OMP_PLACES

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Follow the instructions on the hands-on handout!

Run your benchmark!

- OpenACC
- OpenCL
- OpenMP (2012)