# Hotspots Analysis Intel® VTune TM Profiler

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### Agenda VTune Hotspots

- Overview on this analysis type: views, filters, grouping
- How to run an analysis GUI vs. Command Line
- How to handle MPI programs
- Demo showing typical steps

# Optimize Performance Intel® VTune™ Profiler

#### Get the Right Data to Find Bottlenecks

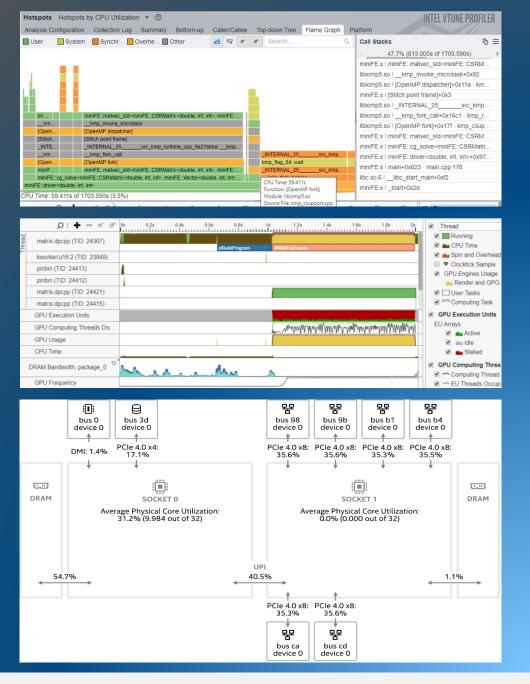
- A suite of profiling for CPU, GPU, FPGA, threading, memory, cache, storage, offload, power...
- Application or system-wide analysis
- DPC++, C, C++, Fortran, Python\*, Go\*, Java\*, or a mix
- Linux, Windows, FreeBSD, Android, Yocto and more
- Containers and VMs

#### Analyze Data Faster

- Collect data HW/SW sampling and tracing w/o recompilation
- See results on your source, in architecture diagrams, as a histogram, on a timeline...
- Filter and organize data to find answers

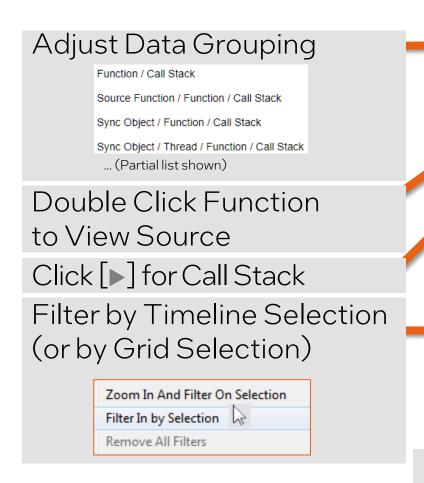
#### Work Your Way

- User interface or command line
- Profile locally and remotely
- GUI (desktop or web) or command line

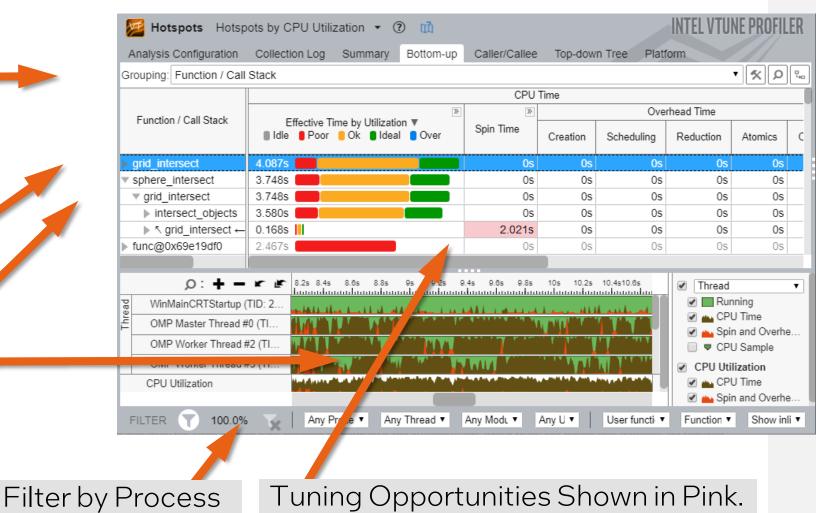


#### Find Answers Fast

Intel® VTune TM Profiler



& Other Controls



## See Profile Data On Source / Asm

Double Click from Grid or Timeline

View Source / Asm or both CPU Time Right click for instruction reference manual Assembly grouping: Address Source Assembly CPU Time: Total Address A Sour CPU Time: T Source Assembly UX4UGCD9 pe ux4uaa49 <Block 53> tmax.x +- taeita.x; U.Z47S lock 44: curpos = nXp; 0x40dcbf 0x40dcbf est esi, esi 0.007s Quick Asm navigation: z 0x40dd07 <Block 50> 0x40dcc1 0.004s 0x40dcc3 Block 45: Select source to highlight Asm mov eax, dword ptr [esi+0x4] 0x40dcc3 0.053s 0.265s 0x40dcc6 mov ecx, dword ptr [edi+0x10] 0.750s cur = g->cells[voxindex]; 0x40dcc9 mov edx, dword ptr [edi+0xc] 0.020s while (cur != NULL) 0.007s 0x40dccc 573 mov eax, dword ptr [eax] 0.055s cmp dword ptr [ecx+eax\*4], edx if (ry->mbox[cur->obj->id] != 0x40dcce 1.177s 574 ry->mbox[cur->obj->id] = ry-> 0.604s 0x40dcd1 jz 0x40dce5 <Block 48> 0.003s 575 0x40dcd3 Block 46: cur->obj->methods->intersect 0.687s 0x40dcd3 mov dword pt [ecx+eax\*4], edx 0.604s 0.423s 0x40dcd6 mov eax, dwo l ptr [esi+0x4] 0.175s cur = cur->next; 578 0x40dcd9 0.005spush edi push eax 0.019s 0x40dcda 0.004s curvox.z += step.z; if (ry->maxdist < tmax.z || curvo 0.011s 0x40dcdb mov ecx, dwor ptr [eax+0x8] 0.027s 0.130s break; 0x40dcde mov eax, dwor ptr [ecx] 0.078s voxindex += step.z\*q->xsize\*q->ys 0x40dce0 call eax

Scroll Bar "Heat Map" is an overview of hot spots

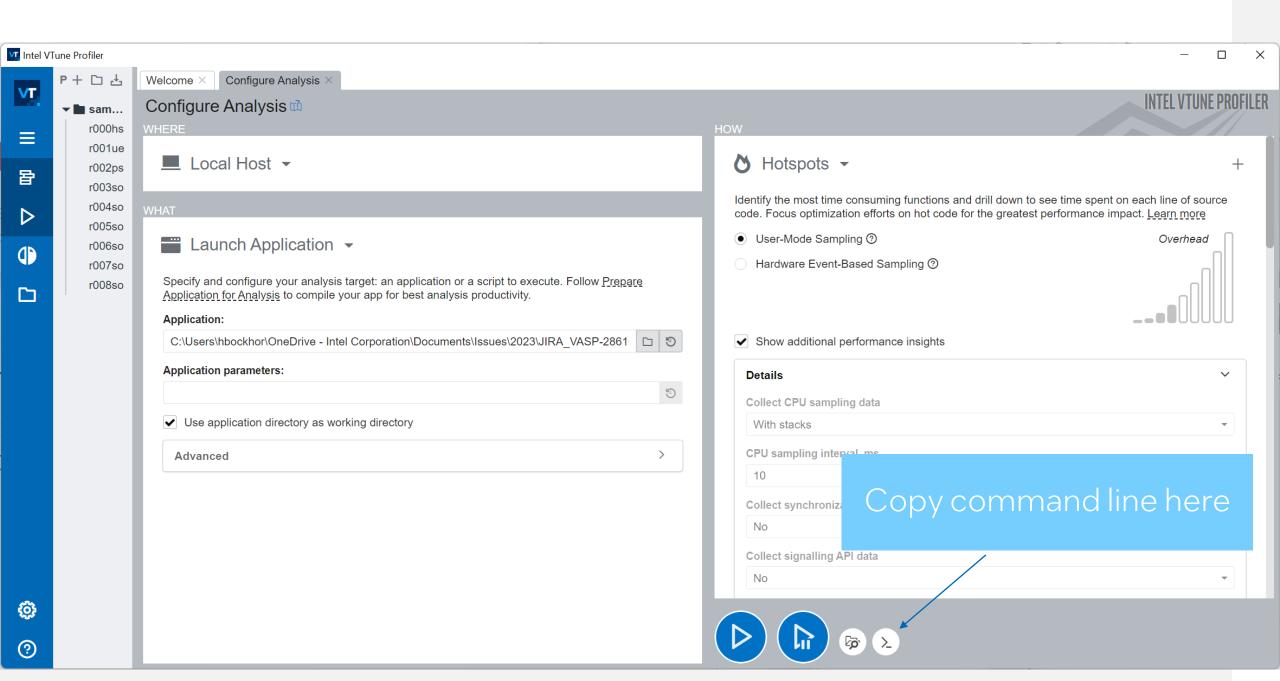
Click jump to scroll Asm

#### Bottom-up tab – most popular tab

- Grouping: different ordering of results –check out different choices
- Source view: double click on function or loop will open another window
   source must be compiled with "-g". Source code must be available
- Zoom and filter in timeline section. Grid will adapt your choice
- Filter process/user code/libraries/loops/system functions (e.g. libc)
- Values in pink: e.g. high overhead like barrier waiting or low cpu utilization.

### How to start an analysis GUI vs CMD

- Start GUI by \$ vtune-gui on command line or by double clicking on Windows
- Click on "Configure Analysis ..."
- 3 Sections: WHERE, WHAT, HOW
- WHERE: local vs remote. We do local here
- WHAT: define your application with parameters and environment
- HOW: Analysis type like "Hotspots" or "APS" with additional parameters



#### Command line

- Copy command line from GUI if you are not sure about the right configuration
- VTune has extensive help menu:
  - \$ vtune -help
  - \$ vtune -help collect
  - \$ vtune -help collect hotspots
- Check "Playbook.txt" for example command lines and additional flags

#### VTune and MPI

- To run VTune in an MPI job you may use the "-gtool" flag
- More convenient is the I\_MPI\_GTOOL environment variable. Example for HPC analysis:

```
$export I_MPI_GTOOL= "vtune -c hpc-performance -r HPC:0"
```

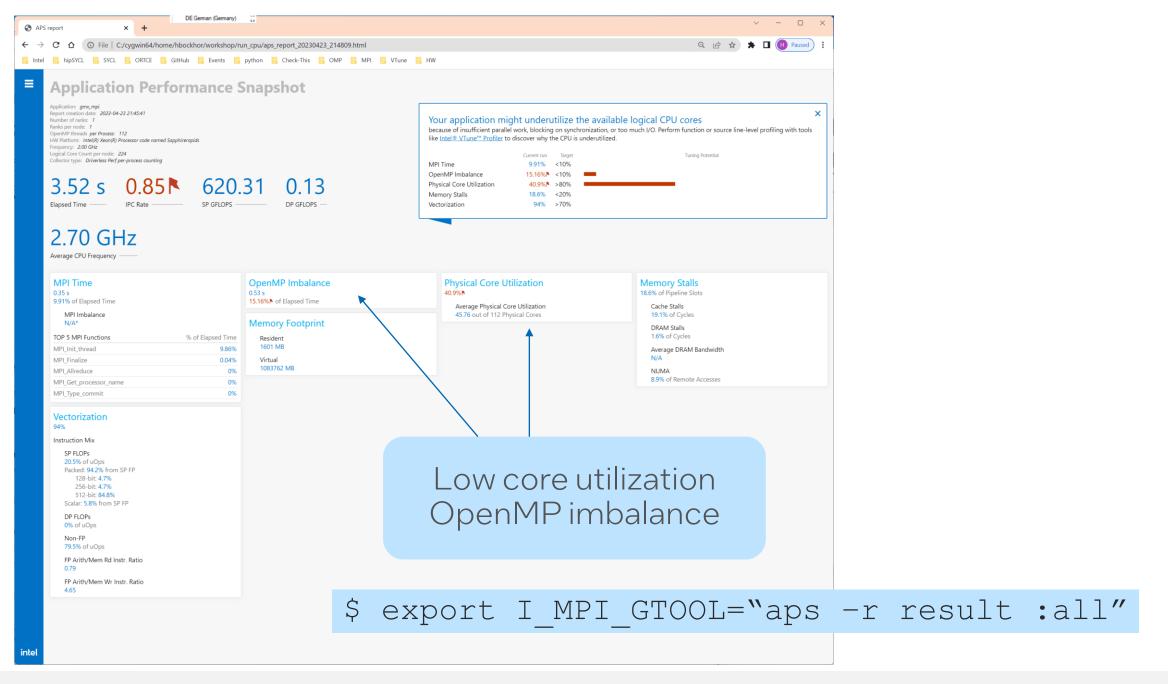
run your program, as usual, under MPI. The setting will collect data on rank #0. Use a list of ranks or : all for multi rank analysis.

More information:

https://www.intel.com/content/www/us/en/develop/documentation/mpi-developer-reference-linux/top/command-reference/mpiexec-hydra/gtool-options.html

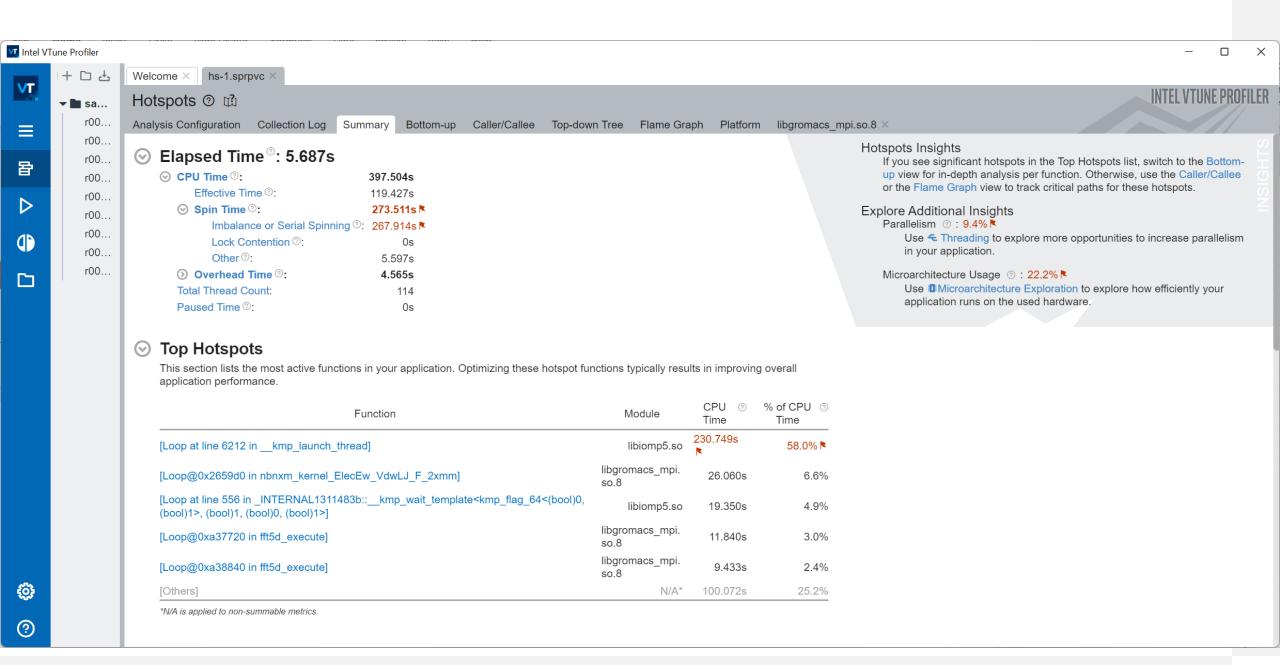
### GROMACS description

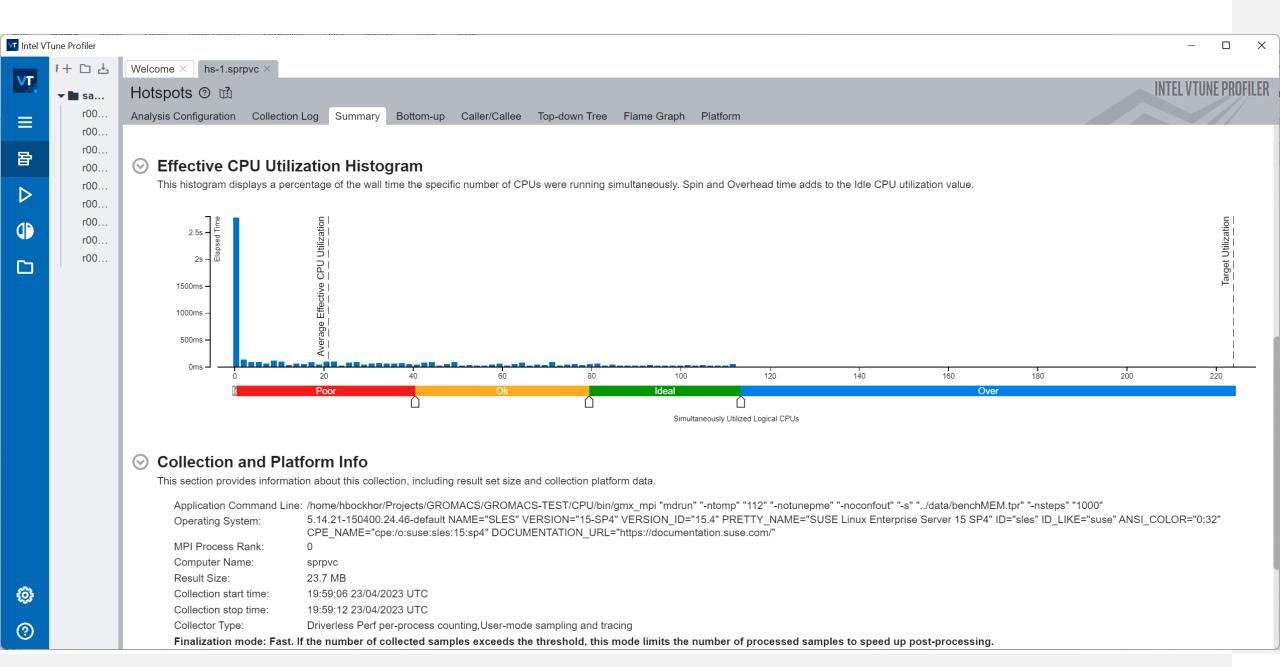
- Molecular Dynamics code with highly tuned CPU and GPU branch
- CPU branch uses AVX-512 intrinsics cmake configuration
- Classical intel compilers (icc,icpc) were deprecated
- Calculation of forces has 2 large parts: direct calculations and large distance forces using FFTs (Particle Mesh Ewald PME)
- Everything is tuned but (intended) naïve execution shows some room for improvement

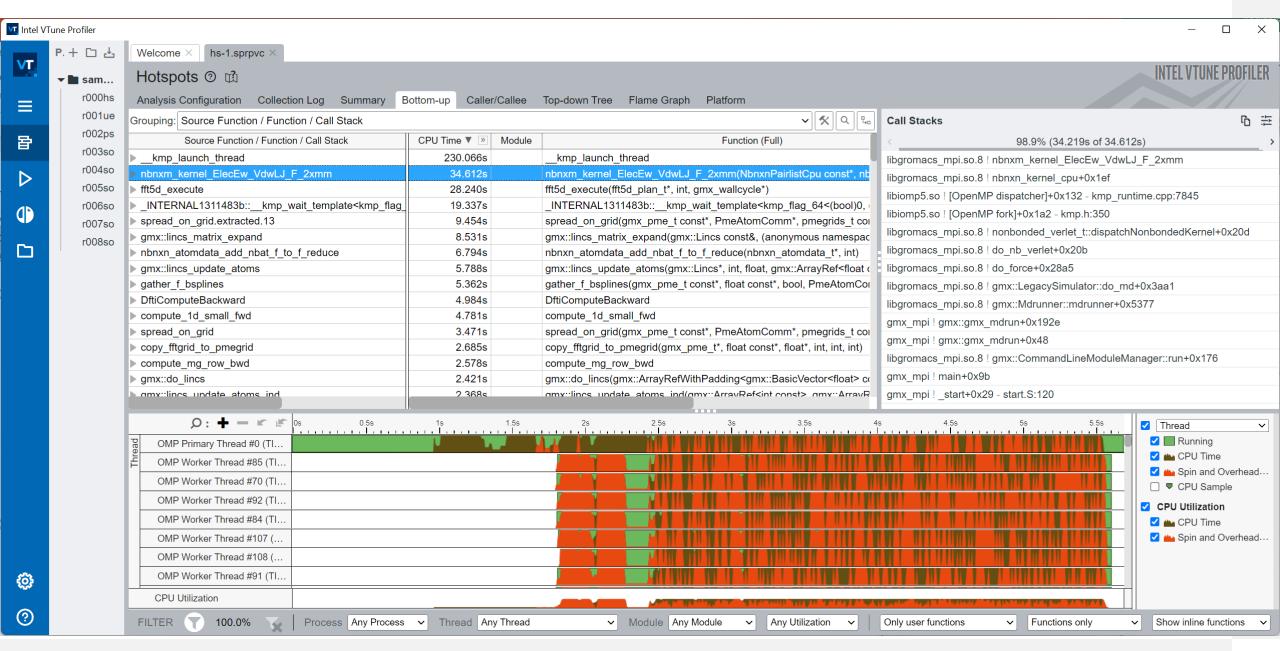


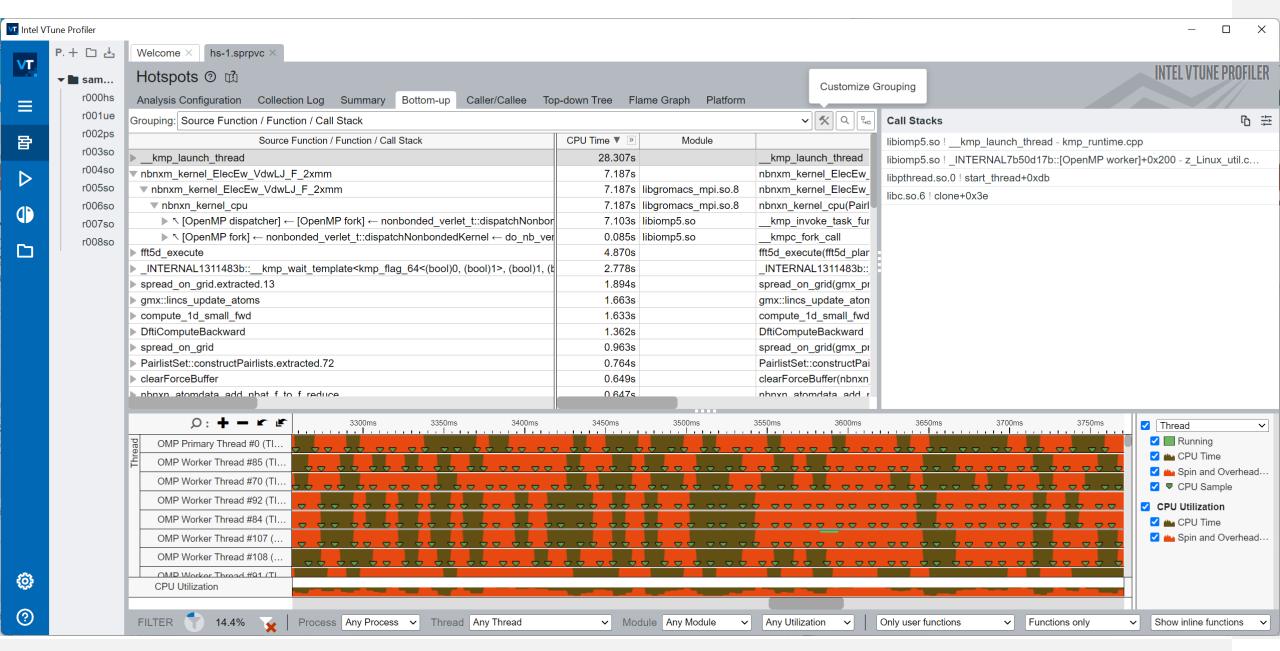
#### APS information

- Low logical core utilization : only single thread per core used
- High OMP imbalance: this can be due to issues with the algorithm or a bad configuration of the environment
- Next step(s): Use VTune with hotspot analysis
- hpc-performance might be better choice, but this comes later ...
- Check command line in the playbook









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