

# Intel Developer Tools and Libraries

The Intel Parallel Studio XE Composer Edition (C/C++/Fortran) is available to all SLAC users. Our license restricts the number of concurrent builds. There are no license restrictions on the runtime libraries.

<https://software.intel.com/en-us/intel-parallel-studio-xe>

We recommend you configure your shell environment by sourcing the appropriate script.

For csh/tcsh, do this:

```
source /afs/slac/package/intel_tools/prod/bin/compilervars.csh intel64
```

For bash-type shells, do:

```
./afs/slac/package/intel_tools/prod/bin/compilervars.sh intel64
```

## OpenMPI with Intel Compilers

We have recompiled a version of the Red Hat OpenMPI package using the Intel compilers. You will need to set your shell environment as described above and also use the **intel-openmpi\_1.8.1-x86\_64** module. Here's a method of using **intel-openmpi\_1.8.1-x86\_64** by default on the bulletmpi cluster. Add the appropriate script from below to your login shell. csh or tcsh users will update .cshrc and bash users will update .bash\_profile or .bashrc.

```
##--FOR CSH or TCSH -----
set bulletcluster = `hostname | grep "^bullet"`
if ($bulletcluster != "") then
source /afs/slac/package/intel_tools/prod/bin/compilervars.csh intel64
eval `/usr/bin/modulecmd csh unload lsf-openmpi_1.5.4-x86_64`
eval `/usr/bin/modulecmd csh load intel-openmpi_1.8.1-x86_64`
endif
#-----

##--FOR BASH -----
bulletcluster=`hostname | grep "^bullet"`
if [ "$bulletcluster" != "" ]; then
. /afs/slac/package/intel_tools/prod/bin/compilervars.sh intel64
eval `/usr/bin/modulecmd sh unload lsf-openmpi_1.5.4-x86_64`
eval `/usr/bin/modulecmd sh load intel-openmpi_1.8.1-x86_64`
fi
#-----
```

[Please see our guidelines for the bulletmpi job queues and OpenMPI](#)