

<b>Software</b>	<b>Link</b>	<b>Description</b>
<b>amber 16.0</b>	<a href="http://ambermd.org/">http://ambermd.org/</a>	Amber is the collective name for a suite of programs that allow users to carry out molecular dynamics simulations
<b>arpack-ng 3.4.0</b>	<a href="https://forge.scilab.org/index.php/p/arpack-ng/">https://forge.scilab.org/index.php/p/arpack-ng/</a>	
<b>bamtools 2.4.1</b>		A programmer's API and an end-user's toolkit for handling BAM files.
<b>blas 3.6.0</b>	<a href="http://www.netlib.org/blas/">http://www.netlib.org/blas/</a>	The Basic Linear Algebra Subprograms library
<b>Cmake 3.5.2</b>	<a href="https://cmake.org/">https://cmake.org/</a>	CMake is a family of tools designed to build, test and package software.
<b>Cmake 2.8.12.2</b>		
<b>comsol 5.2</b>	<a href="https://www.comsol.com/">https://www.comsol.com/</a>	COMSOL Multiphysics® is a general-purpose software platform, based on advanced numerical methods, for modeling and simulating physics-based problems.
<b>dtt 6.1</b>	<a href="http://www.allinea.com/products/dtt">http://www.allinea.com/products/dtt</a>	Parallel debugger supporting a wide range of parallel architectures and models, including MPI, UPC, CUDA and OpenMP.
<b>eigen 3.2.9</b>	<a href="http://eigen.tuxfamily.org/index.php?title=Main_Page">http://eigen.tuxfamily.org/index.php?title=Main_Page</a>	Eigen is a C++ template library for linear algebra: matrices, vectors, numerical solvers, and related algorithms
<b>fastqc 0.11.2</b>	<a href="http://www.bioinformatics.babraham.ac.uk/projects/fastqc/">http://www.bioinformatics.babraham.ac.uk/projects/fastqc/</a>	A quality control tool for high throughput sequence data.
<b>fftw 3.3.5</b>	<a href="http://www.fftw.org/">http://www.fftw.org/</a>	Fast Fourier Transform library
<b>gcc 5.3.1</b>	<a href="https://gcc.gnu.org/">https://gcc.gnu.org/</a>	Various compilers (C, C++, Objective-C, Java, ...)
<b>htslib 1.1.1</b>	<a href="http://www.htslib.org/">http://www.htslib.org/</a>	An implementation of a unified C library for accessing common file formats, such as SAM, CRAM, VCF, and BCF, used for high-throughput sequencing data. It is the core library used by samtools and bcftools.
<b>idl 8.5</b>	<a href="http://www.harrisgeospatial.com/ProductsandSolution">http://www.harrisgeospatial.com/ProductsandSolution</a>	A programming language for extracting visualizations from complex numerical data.

	<a href="s/GeospatialProducts/IDL.aspx">s/GeospatialProducts/IDL.aspx</a>	
<b>impi 5.1.3</b>		
<b>intel 16.0.3</b>		
<b>lammps 20160216</b>	<a href="http://lammps.sandia.gov/">http://lammps.sandia.gov/</a>	LAMMPS Molecular Dynamics Simulator
<b>lapack 3.6.1</b>	<a href="http://www.netlib.org/lapack/">http://www.netlib.org/lapack/</a>	Numerical linear algebra package libraries
<b>Mathematica</b>	<a href="https://www.wolfram.com/mathematica/">https://www.wolfram.com/mathematica/</a>	
<b>matlab R2016a</b>	<a href="http://www.mathworks.com">http://www.mathworks.com</a>	MATLAB(R) is a high-level language and interactive environment for numerical computation, visualization, and programming.
<b>metis 5.1.0</b>	<a href="http://glaros.dtc.umn.edu/gkhome/metis/metis/overview">http://glaros.dtc.umn.edu/gkhome/metis/metis/overview</a>	Serial Graph Partitioning and Fill-reducing Matrix Ordering
<b>mpich 3.2</b>	<a href="https://www.mpich.org/">https://www.mpich.org/</a>	A high-performance implementation of MPI
<b>mpir 1.3.1</b>	<a href="http://mpir.org/">http://mpir.org/</a>	A library for arbitrary precision arithmetic
<b>netcdf 4.4.1</b>	<a href="http://www.unidata.ucar.edu/software/netcdf/">http://www.unidata.ucar.edu/software/netcdf/</a>	Libraries for the Unidata network Common Data Form
<b>netcdf-fortran 4.4.4</b>		
<b>octave 3.8.2</b>	<a href="https://www.gnu.org/software/octave/">https://www.gnu.org/software/octave/</a>	GNU Octave is a high-level interpreted language, primarily intended for numerical computations.
<b>openmpi 1.10.1</b>	<a href="https://www.open-mpi.org/">https://www.open-mpi.org/</a>	A powerful implementation of MPI/SHMEM
<b>petsc 3.7</b>	<a href="https://www.mcs.anl.gov/petsc/">https://www.mcs.anl.gov/petsc/</a>	Portable, Extensible Toolkit for Scientific Computation
<b>pgi 16.4</b>	<a href="https://www.pgroup.com/about/">https://www.pgroup.com/about/</a>	Portland Group (PGI) is a premier supplier of software compilers and tool for parallel computing, known as PGI products. The Portland Group offers optimizing parallel FORTRAN 2003, C99, and C++ compilers and tools for workstations, servers and clusters running Linux, MacOS, and Windows
<b>pointwise</b>	<a href="http://www.pointwise.com/">http://www.pointwise.com/</a>	

<b>python 3.3.2</b>	<a href="https://www.python.org/">https://www.python.org/</a>	
<b>python 2.7.8</b>		
<b>python 2.6.6</b>		
<b>python-astropy</b>		
<b>python-ipython 0.13.2</b>		
<b>python-matplotlib 0.99.1.2</b>		
<b>python-mne 0.12.0</b>		
<b>python-numdisplay 1.5.6</b>		
<b>python-numpy 1.4.1</b>		
<b>python-scipy 0.7.2</b>		
<b>quantum-espresso 5.4.0</b>	<a href="http://www.quantum-espresso.org/">http://www.quantum-espresso.org/</a>	An integrated suite of Open-Source computer codes for electronic-structure calculations and materials modeling at the nanoscale
<b>R 3.3.1</b>	<a href="https://www.r-project.org/">https://www.r-project.org/</a>	A language for data analysis and graphics
<b>r-iterators 1.0.8</b>		
<b>r-rlecuyer 0.3-4</b>		
<b>r-rmpi 6.0-6</b>		
<b>r-snow 0.4-2</b>		
<b>totalview 8.15.10</b>	<a href="http://www.roguewave.com/products-services/totalview">http://www.roguewave.com/products-services/totalview</a>	Parallel debugger supporting a wide range of parallel architectures and models, including MPI, UPC, CUDA and OpenMP.
<b>trinity 2.2.0</b>	<a href="https://github.com/trinityrnaseq/trinityrnaseq/wiki">https://github.com/trinityrnaseq/trinityrnaseq/wiki</a>	Tools for processing RNA-seq data
<b>udunits 2.2.20</b>	<a href="http://www.unidata.ucar.edu/software/udunits/">http://www.unidata.ucar.edu/software/udunits/</a>	A library for manipulating units of physical quantities
<b>vasp 5.4.1</b>	<a href="https://www.vasp.at/">https://www.vasp.at/</a>	Program for atomic-scale materials modeling
<b>xblas 1.0.248</b>	<a href="http://www.netlib.org/xblas/">http://www.netlib.org/xblas/</a>	Extra Precise Basic Linear Algebra Subroutines
<b>xz 5.2.2</b>	<a href="http://tukaani.org/xz/">http://tukaani.org/xz/</a>	LZMA compression utilities