	A	S	Т
1	Package Name	Software Description	Project URL
2	Autoconf	An extensible package of M4 macros that produce shell scripts to automatically configure software source code packages	https://www.gnu.org/software/autoconf/
3	Automake		www.gnu.org/software/automake
4	Libtool		www.gnu.org/software/libtool
5	bamtools	BamTools: a C++ API for reading/writing BAM files.	https://github.com/pezmaster31/bamtools
6	Biopython (Python module)	Biopython is a set of freely available tools for biological computation written in Python by an international team of developers	www.biopython.org/
7	blas	The BLAS (Basic Linear Algebra Subprograms) are routines that provide standard building blocks for performing basic vector and matrix operations.	http://www.netlib.org/blas/
8	boost	Boost provides tree peer-reviewed portable C++ source libraries.	http://www.boost.org
9	Civiake	Cross-platform, open-source build system. Crivate is a ramily or tools designed to build, test and package software	nttp://www.cmake.org/
10	Cython (Python module)	The Cython compiler for writing C extensions for the Python language	nttps://www.pytnon.org/
	Doxygen	FEmpera is the leading multimedia framework able to decode encode transcode mux demux stream filter and play pretty much anything that humans	mp://www.doxygen.org/
		and machines have created it supports the most obscure ancient formats up to the cuttion edge. No matter if they were designed by some standards	
12	ffmpeg	committee the community or a cororation.	https://www.ffmpeq.org
	1 - 22	FFTW is a C subroutine library for computing the discrete Fourier transform (DFT) in one or more dimensions, of arbitrary input size, and of both real and	
13	fftw	complex data (as well as of even/odd data, i.e. the discrete cosine/sine transforms or DCT/DST).	http://www.fftw.org/download.html
14	git	Git is a distributed version control system designed to handle everything from small to very large projects with speed and efficiency.	http://git-scm.com/
		GMP is a free library for arbitrary precision arithmetic, operating on signed integers, rational numbers, and floating-point numbers. There is no practical	
		limit to the precision except the ones implied by the available memory in the machine GMP runs on. GMP has a rich set of functions, and the functions	
15	gmp	have a regular interface.	https://gmplib.org
		The GNU Compiler Collection includes front ends for C, C++, Objective-C, Fortran, Java, Ada, and Go, as well as libraries for these languages (libstdc++,	
16	GNU Compiler Collection (GCC)		nttps://gcc.gnu.org/
17	CNULEmana	UNU Emacs is an extensible, customizable text editor—and more. At its core is an interpreter for Emacs Lisp, a dialect of theLisp programming language	www.gev.co/opticipe/oppop/
17	GINU ETHAGS	with extensions to support BXI Building. Counted to a particular command line driven graphing utility for Linux, OS/2, MS Mindows, OSX, VMS, and many other statistics. The source and a in	www.gnd.org/sortwate/efflats/
18	anuplot	convioluted but freely distributed (i.e. you don't have to pay for it)	http://www.apuplet.info
19	Graphviz	Graph Visualization Software	http://www.graphviz.org/
20	hdf5	HDF5 (Hierarchical Data Format version 5) from HDFGroup	http://www.hdfaroup.org/HDF5/
		ImageMagick® is a software suite to create, edit, compose, or convert bitmap images. It can read and write images in a variety of formats (over 100)	
		including DPX, EXR, GIF, JPEG, JPEG-2000, PDF, PNG, Postscript, SVG, and TIFF. Use ImageMagick to resize, flip, mirror, rotate, distort, shear and	
21	imagemagick	transform images, adjust image colors, apply various special effects, or draw text, lines, polygons, ellipses and Bézier curves.	http://imagemagick.org
22	isl	ISL (Integer Set Library) is a library for manipulating sets and relations of integer points bounded by linear constraints.	http://isl.gforge.inria.fr
		Java is a general-purpose, concurrent, class-based, object-oriented language that is specifically designed to have as few implementation dependencies as	
23	Java	possible.	http://java.com/
		Julia is a high-level, high-performance dynamic programming language for technical computing, with syntax that is familiar to users of other technical	
24	julia	computing environments.	http://julialang.org
25	lanask	LAPACK is written in Fortran 90 and provides routines for solving systems of simultaneous linear equations, least-squares solutions of linear systems of gruntices elicencilus architecture and elegication under sublighted and the second sec	http://www.petlik.org/opeel/
23	араск	equations, eigenvalue problems, and singular value problems. Manuruli, is a cross platform, distributed revision explicit for software developers. It is maintuimplemented using the Puthen programming language	nttp://www.netito.org/napack/
26	mecurial	mercular is a closs-plationin, distributed revision control to no software developers, it is mainly implemented using the Fython programming language, but includes a binary diffi implementation written in C	http://en.wikipedia.org/wiki/Mercurial
20	mecunai	Durincudes a Dinary dim implementation written in C.	nto.//en.wikipedia.org/wiki/metcunar
		the IEEE-754 standard for fixed precision real floating point numbers to complex numbers, providing well-defined semantics for every operation. At the	
27	mpc	same time, speed of operation at high precision is a maior design goal.	http://www.multiprecision.org/index.php?prog=mpc
		The MPFR library is a C library for multiple-precision floating-point computations with correct rounding. MPFR has continuously been supported by the	
		INRIA and the current main authors come from the Caramel and AriC project-teams at Loria (Nancy, France) and LIP (Lyon, France) respectively; see	
28	mpfr	more on the credit page. MPFR is based on the GMP multiple-precision library.	http://www.mpfr.org
29	mpich	MPICH is a high-performance and widely portable implementation of the Message Passing Interface (MPI) standard (MPI-1, MPI-2 and MPI-3).	http://www.mpich.org
30	nose (Python module)	Module for unit testing in Python	https://github.com/nose-devs/nose/
		GNU Octave is a high-level interpreted language, primarily intended for numerical computations. It provides capabilities for the numerical solution of linear interpreter and explore and for an effort of the numerical burnelistic thread of the numerical solution of linear	
31	octave	and nonlinear problems, and for performing other numerical experiments.	https://www.gnu.org/software/octave/
20	ananmai	The upper Miri Project is an open source Message Passing Interface implementation that is developed and maintained by a consortium of academic,	
32	openinpi	research, and industry partners.	http://www.open-htp.org
34	python3 (+ all the available modules)	Tyton is an object oriented, interpreted, and interactive programming language. (version 2)	https://www.putnor.org/downlads/releases/putnor.342/
35	root	ROOT, an OO framework for large-scale scientific data analysis and data mining	http://root.cem.ch/drupal/
36	ruby (+ all the available modules)	A dynamic, open source programming language with a focus on simplicity and productivity.	https://www.ruby-lang.org/en/
		Suite of sparse matrix algorithms. Includes UMFPACK, CHOLMOD, SPQR, ordering methods (AMD, CAMD, COLAMD, and CCOLAMD) and more. Can do	
37	SuiteSparse	QR factorization on the GPU	http://faculty.cse.tamu.edu/davis/suitesparse.html
38	torque	Loads the Torque batch computing environment	http://www.adaptivecomputing.com/
39	valgrind	Valgrind is an instrumentation framework for building dynamic analysis tools.	http://valgrind.org
	-176	zito is designed to be a tree, general-purpose, legally unencumbered that is, not covered by any patents lossless data-compression library for use on	
40	ZIID	virtually any computer naroware and operating system.	nttp://www.zilio.net
41	asl	Amine DD i is an auvanceu deudyging tool deel no scalat, multi-tineadeu, a to arge-scale paratier apprications. The GNU Scientific Linsarv (GSL) is a numerical library for C and C++ programmers.	http://www.aminea.com/products/downloads/gel-support
43	h5utils	Stutis is a set of utilities for visualization and conversion of scientific data in the free portable HDE5 format	http://www.gito.org/software.ga/
		Hypre is a massively parallelizable linear solver library written and designed for use in C++. It provides many nice features and objects beyond simple	
44	hypre	"matrix" or "vector" objects - these include grids, meshes, and stencils. It was developed at Lawrence Livermore National Lab.	http://charlesmartinreid.com/wiki/Hypre, http://acts.nersc.gov/hypre/
		Intel® Parallel Studio XE 2015 tool suite simplifies the design, development, debug, and tuning of code that utilizes parallel processing to boost application	
45	intel	performance. Get more application performance with less effort on compatible Intel® processors and coprocessors.	https://software.intel.com/en-us/intel-compilers, https://software.intel.com/en-us/intel-parallel-studio-xe
		METIS is a set of serial programs for partitioning graphs, partitioning finite element meshes, and producing fill reducing orderings for sparse matrices. The	
		algorithms implemented in METIS are based on the multilevel recursive-bisection, multilevel k-way, and multi-constraint partitioning schemes developed in	
46	metis		http://glaros.dtc.umn.edu/gkhome/views/metis
47	mki mniak 2	Intel Warn Remei Library MICH2 is a biote active second widely particular particular of the Massage Dessing Interface (MDI) standard (MDI 4, MDI 4, STAL MDI 4).	nttps://sortware.intel.com/en-us/intel-mki
48	11ptch2	wmmunizes a migreperiorinance and widely portable implementation or the message Passing Interface (MPI) standard (MPI-1, MPI-2 and MPI-3). The netCDE Operator (MCO) tookit manipulates and analyzes data stored in perCDE concercible formate.	http://www.mpicn.org
49 50	netodf	Interfaces operates (recy) toolkit manipulates and analyzes data stored in netCDP+dCdessible formats NetCDE version 4.3.2 (with NetCDE Fortran I brazies)	http://www.unidata.ucar.edu/software/oetcdf/
50		DoenCV (Doen Source Computer Vision Librario) OpenCV (Doen Source Computer Vision Librario) is an open source computer vision and machine learning software library. OpenCV was hull to provide a	nings// mmm.uningsits.oodi.codi/softmarc/ifotodi/
51	opency	common infrastructure for computer vision applications and to accelerate the use of machine percention in the commercial products	http://opency.org
- · ·		ParMETIS is an MPI-based parallel library that implements a variety of algorithms for partitioning unstructured graphs, meshes, and for computing fill-	
52	parmetis	reducing orderings of sparse matrices.	http://glaros.dtc.umn.edu/gkhome/metis/parmetis/overview
		PETSc, pronounced PET-see (the S is silent), is a suite of data structures and routines for the scalable (parallel) solution of scientific applications modeled	
53	petsc	by partial differential equations.	http://www.mcs.anl.gov/petsc/

	Δ	S. S	Т
54	pgi	Portiand Group Compiler Accelerator Suite (c, c++ and ronran Server x86_64)	https://www.pdc.kth.se/software/compilers/pgi/general.html
55	scalapack	ScaLAPACK is a library of high-performance linear algebra routines for parallel distributed memory machines.	http://www.netlib.org/scalapack/
		SLEPc is a software library for the solution of large scale sparse eigenvalue problems on parallel computers. It is an extension of PETSc and can be used	
56	slepc	for linear eigenvalue problems in either standard or generalized form, with real or complex arithmetic.	http://www.grycap.upv.es/slepc/
		SUNDIALS was implemented with the goal of providing robust time integrators and nonlinear solvers that can easily be incorporated into existing	
57	sundials	simulation codes.	http://computation.llnl.gov/casc/sundials/main.html
		Total View is a GLIL-based source code debugger and defect analysis tool that provides unprecedented control over processes and thread execution and	
59	totalviow	visibility into program that and variable.	http://www.roguowovo.com/oroducts/totalview.ocpv
30	totalview	visionity into program state and variables.	http://www.logdewave.com/products/totamew.aspx
	and and the	The oboint's package supports units of physical quantities, its C library provides for antimetic manipulation of units and for conversion of numeric values	
59	udunits	between compatible units.	nttp://www.unidata.ucar.edu/sontware/udunits/
60	virtualenv	virtualenv is a tool to create isolated Python environments.	http://virtualenv.readthedocs.org/en/latest/virtualenv.html
61	vtk	The Visualization Toolkit (VTK) is an open-source, freely available software system for 3D computer graphics, image processing and visualization.	http://www.vtk.org/VTK/resources/software.html
62	xblas	XBLAS - Extra Precise Basic Linear Algebra Subroutines	http://www.netlib.org/xblas
		Fast C++ matrix library with easy to use functions and syntax, deliberately similar to Matlab. Uses template meta-programming techniques. Also provides	
63	armadillo	efficient wrappers for LAPACK, BLAS and ATLAS libraries, including high-performance versions such as Intel MKL, AMD ACML and OpenBLAS.	http://arma.sourceforge.net
64	bdwgc	The Boehm-Demers-Weiser conservative C/C++ Garbage Collector (libgc, bdwgc, boehmgc)	http://www.hboehm.info/gc/
		The reference BLAS is a freely-available software package. It is available from netlib via anonymous ftp and the World Wide Web. Thus, it can be	
65	cblas	included in commercial software packages (and has been). We only ask that proper credit be given to the authors.	http://www.netlib.org/blas/#_cblas.http://www.prism.gatech.edu/~ndantam3/cblas-doc/doc/html/main.html
		C on G is a free software and library to generate onde for scanning Z-polyhedra. That is it finds a code (e.g. in C EORTRAN) that reaches each integral	
66	cloog	Debet is a new solvane and industrial could be solvaning 2 polynoida. That is, it indus a could (e.g. in 0, it or there w) that reaches each integration of the solution of	http://www.cloog.org
00	cibog	point of one parameterized polyheara.	http://www.clobg.org
		The twining of the control of the co	
		CODA rookit includes a complier for rivible GPOs, main libraries, and tools for debugging and optimizing the performance of your applications. Four	
	0100 T 113	also into programming guides, user manuals, API reference, and other documentation to help you get started quickly accelerating your application with	
67	CUDA Toolkit	GPUS.	nttps://developer.nvidia.com/cuda-toolkit
68	deal.II	A C++ software library supporting the creation of finite element codes and an open community of users and developers.	http://www.dealii.org
		The aim of the decorator module it to simplify the usage of decorators for the average programmer, and to popularize decorators by showing various non-	
69	Decorator (Python module)	trivial examples	https://pypi.python.org/pypi/decorator
70	dx	OpenDX (IBM Visualization Data Explorer) Data Explorer is a system of tools and user interfaces for visualizing data.	http://www.opendx.org/highlights.php?highlight=inaction/tutorials/
		Eclipse is famous for our Java Integrated Development Environment (IDE), but our C/C++ IDE and PHP IDE are pretty cool too. You can easily combine	
		language support and other features into any of our default packages and the Eclinse Marketolace allows for virtually unlimited customization and	
71	Eclipse	extension.	http://www.eclipse.org/users/
	LCIIpse	Extension.	http://www.eclipse.org/users/
		I fullie is a distributed, reliable, and available service for enciencing, augregating, and moving large aniouns of log data. It has a simple and	
-		nexible a childcure based on streaming data nows, it is robust and radit tolerant with tunable reliability mechanisms and many failover and recovery	
72	flume	mechanisms. It uses a simple extensible data model that allows for online analytic application.	nttps://flume.apache.org/
-		GDB, the GNU Project debugger, allows you to see what is going on inside another program while it executes or what another program was doing at	
73	gab		nttp://www.gnu.org/software/gdb/
		GMT is a collection of public-domain Unix tools that allows you to manipulate x,y and x,y,z data sets (filtering, trend fitting, gridding, projecting, etc.) and	
		produce PostScript illustrations ranging from simple x-y plots, via contour maps, to artificially illuminated surfaces and 3-D perspective views in black/white	
		or full color. Linear, log10, and power scaling is supported in addition to 25 common map projections. The processing and display routines within GMT are	
74	gmt	completely general and will handle any (x,y) or (x,y,z) data as input.	http://tossies.org/linux/misc/GMT/gmt-4.5.12-src.tar.gz/gmt-4.5.12/src/GMT.1
		Grace is a WYSIWYG 2D plotting tool for the X Window System and M*tif. Grace runs on practically any version of Unix-like OS. As well, it has been	
75	grace	successfully ported to VMS, OS/2, and Win9*/NT/2000/XP (some minor functionality may be missing, though).	http://plasma-gate.weizmann.ac.il/Grace/
		GStreamer is a library for constructing graphs of media-handling components. The applications it supports range from simple Ogg/Vorbis playback,	
76	gstreamer	audio/video streaming to complex audio (mixing) and video (non-linear editing) processing.	http://gstreamer.freedesktop.org
		Guile is the GNU Ubiquitous Intelligent Language for Extensions, the official extension language for the GNU operating system. Guile is a library designed	
		to help programmers create flexible applications. Using Guile in an application allows the application's functionality to be extended by users or other	
77	guile	programmers with plug-ins, modules, or scripts.	https://www.gnu.org/software/guile/download.html
		Harminv is a free program (and accompanying library) to solve the problem of harmonic inversion — given a discrete-time, finite-length signal that consists	
		of a sum of finitely-many sinusoids (possibly exponentially decaying) in a given bandwidth, it determines the frequencies, decay constants, amplitudes,	
78	harminv	and phases of those sinusoids.	http://ab-initio.mit.edu/wiki/index.php/Harminv
79	hdf4	HDF4 (Hierarchical Data Format version 4) from HDFGroup	http://www.hdfgroup.org/products/hdf4/
		ITK is an open-source software toolkit for performing registration and segmentation. Segmentation is the process of identifying and classifying data found	
80	InsightToolkit	in a digitally sampled representation.	http://www.itk.org/ITK/project/project.html
81	jags	JAGS is Just Another Gibbs Sampler. It is a program for analysis of Bayesian hierarchical models using Markov Chain Monte Carlo (MCMC) simulation	http://mcmc-jags.sourceforge.net
82	libctl	ibctl is a free Guile-based library implementing flexible control files for scientific simulations.	http://ab-initio.mit.edu/wiki/index.php/Libctl
		A Portable Foreign Function Interface Library - The libfif library provides a portable, high level programming interface to various calling conventions. This	
83	libffi	allows a programmer to call any function specified by a call interface description at run-time.	https://sourceware.org/libffi/
84	libsbml	LibSBML is a free, open-source programming library to help you read, write, manipulate, translate, and validate SBML files and data streams.	http://sbml.org/Software/libSBML
85	libunistring	libunistring - This library provides functions for manipulating Unicode strings and for manipulating C strings according to the Unicode standard.	https://www.gnu.org/software/libunistring/
		Multiple Precision Integers and Rationals (MPIR) is an open-source software multiprecision integer library forked from the GNU Multiple Precision	
86	mpir	Arithmetic Library (GMP) project. It consists of much code from past GMP releases, and much newer original contributed code	http://mpir.org
		The NASMX Project is a collection of macros, includes and examples, which are meant to help NASM programmers develop 32-bit and 64-bit BSD. Linux.	
87	namsxz	Windows, and XBOX software in a fraction of the time it normally takes.	http://www.nasm.us
		The p4est software library enables the dynamic management of a collection of adaptive octrees, conveniently called a forest of octrees. p4est is designed	
88	p4est	to work in parallel and scales to hundreds of thousands of processor cores.	http://www.p4est.org
		PAPI - (Performance Application Programming Interface) provides the tool designer and application engineer with a consistent interface and methodology	
89	papi	for use of the performance counter hardware found in most major microprocessors.	http://icl.cs.utk.edu/papi/
		Program Database Toolkit (PDT) is a framework for analyzing source code written in several programming languages and for making rich program	· · · · ·
90	pdt	knowledge accessible to developers of static and dynamic analysis tools.	http://www.cs.uoregon.edu/research/odt/home.php
		The Parma Polyhedra Library (PPL) provides numerical abstractions especially targeted at applications in the field of analysis and verification of complex	
91	laa	systems.	http://buasena.com/products/pp/
92	sac	SAC (Seismic Analysis Code) is a general purpose interactive program designed for the study of sequential signals, especially time series data	https://seiscode.iris.washington.edu/projects/sac
~~		Score is an Open Source software construction tool—that is a next-openeration build tool. Think of Score as an improved cross-relation substitute for	
		the classic Make utility with integrated functionality similar to autoconfoutomed and complete caches used as manyority of the classic Make utility with integrated functionality similar to autoconfoutomed and complete caches used as cache in short SCone is an easier more	
02	scons	The backs may may may may be an any may be an any and a set of the	http://www.scops.org
33	30013	SPDARKS is a parallel Monte Carlo code for on lattice and off-lattice models that includes algorithms for kinotic Monte Carlo (KMC), reliables literation	Intp.//www.soons.org
04	sonarke	Anna Carlo (KMC), and Matronolis Monte Carlo (MMC)	http://coport/c condia.gov
94	σμαινο	mone can entword, and metropolis mone can entwork.	III(µ.//sµµaitossaitula.\$UV
05	outo	owners a sonware development toor that connects programs written in C and C++ with a variety or nigh-level programming languages. SWIG is used with different home of target longuages including common containing longuages such as a longuage. Det Det Det as a co	http://www.evid.org
95	swig	umeren types on target languages including common scripting languages such as Javaschipt, Peri, PHP, Pytron, Tci and Ruby.	ntp://www.swig.org
		The renormance system with a portable proming and tracing tookit for performance analysis of parallel programs written in Fortran, C, C++, UPC, Java,	
96	tau	Pyrion.	http://www.cs.uoregon.edu/research/tau/home.php
		I mux terminal multiplexer- lets you switch easily between several programs in one terminal, detach them (they keep running in the background) and	
97	tmux	reattach them to a different terminal.	http://tmux.sourcetorge.net

	Α	S	Т
	8	0	I
98	trinity	Irinity is an integrated development/debugging environment (IDE) for Derric, a domain-specific language to declaratively describe binary file formats.	https://github.com/jvdb/trinity
99	visit	Visit is an Open Source, interactive, scalable, visualization, animation and analysis tool.	https://wci.llnl.gov/simulation/computer-codes/visit
		matplotlib is a python 2D plotting library which produces publication quality figures in a variety of hardcopy formats and interactive environments across	
100	matplotlib (Python module)	platforms	http://matplotlib.org/
		Python library that provides a multidimensional array object, various derived objects (such as masked arrays and matrices), and an assortment of routines	
		for fact operations on arrays including mathematical logical shape manipulation sorting selecting 1/O discrete Fourier transforms basic linear algebra	
101	NumPy (Python modulo)	basic statistical portations of analys, motioning mationalizati, logical, shape manipulation, soluting, solecuting, i/o, discrete roundin transforms, basic finear algebra, basic statistical portations mandom simulation.	http://www.pumpy.org/
101	Numey (Eython module)		ntp://www.numpy.org/
		SCIPy contains additional routines needed in scientific work: for example, routines for computing integrals numerically, solving differential equations,	
102	SciPy (Python module)	optimization, and sparse matrices	www.scipy.org/
103	cdo	Climate Data Operators (CDO) - Command line Operators for Climate and NWP model Data.	https://code.zmaw.de/projects/cdo
104	gdal	GDAL - Geospatial Data Abstraction Library - a translator library for raster and vector geospatial data formats	http://www.gdal.org
		GEOS (Geometry Engine - Open Source) is a C++ port of the Java Topology Suite (JTS). As such, it aims to contain the complete functionality of JTS in	
		C++. This includes all the OpenGIS Simple Features for SQL spatial predicate functions and spatial operators, as well as specific JTS enhanced topology	
105	deos	functions.	http://trac.osgeo.org/geos/
	3	The Grid Analysis and Disnlay System (GrADS) is an interactive desitor tool that is used for easy access, manipulation, and visualization of earth science	······································
106	arada	The one Analysis and Display Gystein (ChADG) is an interactive deskep too that is deed for easy access, manipulation, and visualization of caraf science	http://iaos.org/arads/
100	graus	uala. Di la la fanta de la sistifica en energia de serve de serve dissidirente entre fanta de la construction de serve	nup//iges.org/glads/
107	101	IDE is the trusted scientific programming language used across disciplines to extract meaningful visualizations out or complex numerical data.	nttp://exellsvis.com/Home.aspx
108	nci	NCL - NCAR Command Language - Meteorological Graphics Plotting System	nttp://www.nci.ucar.edu
109	ncview	Noview is a visual browser for netCDF format files.	http://meteora.ucsd.edu/~pierce/ncview home page.html
110	panoply	Panoply is a cross-platform application that plots geo-gridded and other arrays from netCDF, HDF, GRIB, and other datasets.	http://www.giss.nasa.gov/tools/panoply/
111	proj	PROJ.4 - Cartographic Projections Library	http://trac.osgeo.org/proj/
1		Delft3D is a world leading 3D modeling suite to investigate hydrodynamics, sediment transport and morphology and water quality for fluvial estuaring and	
112	delft3d	coastal environments. As per 1 January 2011, the Delfab flow (FLOW), morphology (MOR) and waves (WAVE) modules are available in open source	http://oss.deltares.pl/web/delft3d
112	enrnt	a river dynamics simulator	https://doi.org/10.100/100100
113	opint	a more oprianilized annotation.	nteps/gende.commany-ne/or/twn
		minute refers to two unings, a set or indirectian metrianical rorce neitos for the simulation or bioindirecties (which are in the public domain, and are used in a	http://www.hours.down
114	amper	variety or simulation programs); and a package or molecular simulation programs which includes source code and demos.	nttp://ambermd.org
1		peditools utilities are a swiss-army knife of tools for a wide-range of Life Sciences analysis tasks. The most widely-used tools enable genome arithmetic:	
115	bedtools	that is, set theory on the genome.	http://bedtools.readthedocs.org/en/latest/
		The Basic Local Alignment Search Tool (BLAST) finds regions of local similarity between sequences. The program compares nucleotide or	
		proteinsequences to sequence databases and calculates the statistical significance of matches. BLAST can be used to infer functional and evolutionary	
116	blast	relationships between sequences as well as help identify members of gene families	http://blast.pcbi.plm.pib.gov/Rlast.cgi
110	biast	Power on ultrafact memory officiant chart and aligner that are on our DNA concept (reade) to the human genome at a rate of over 25 million 25 ha	nto //ordschob/ninninnigov/bidscogr
447	h an dia	bowie is an uitalast, memory-encient short read aligner. It aligns short DrvA sequences (reads) to the numan genome at a rate of over 25 million 55-bp	
117	bowtie		nttp://dowtie-bio.sourcerorge.nevmanuai.sntmi
		BWA is a program for aligning sequencing reads against a large reference genome (e.g. human genome). It has two major components, one for read	
118	bwa	shorter than 150bp and the other for longer reads.	http://bio-bwa.sourceforge.net
		Cufflinks assembles transcripts, estimates their abundances, and tests for differential expression and regulation in RNA-Seq samples. It accepts aligned	
		RNA-Seq reads and assembles the alignments into a parsimonious set of transcripts. Cufflinks then estimates the relative abundances of these transcripts	
119	cufflinks	based on how many reads support each one, taking into account biases in library preparation protocols.	http://cufflinks.cbcb.umd.edu
120	express	express is a streaming tool for quantifying the abundances of a set of target sequences from sampled subsequences.	http://bio.math.berkelev.edu/eXpress/
		The FASTA programs can be used to search protein and DNA sequence databases and to confirm the statistical significance of a match by comparing	
121	facto	The align programs can be added to seen produced by shuffled conjugates, and to commit the statistical significance of a match by comparing	http://foculty.virginia.odu/urpogrop/focto/
121	factor	and angiment source of a distribution of scores produced by shared sequences.	http://naouty.wigina.cou/wipca.sou/asia/
122	lasiqu	A quanty control tool for high through put sequence data	http://www.bioinformatics.babranam.ac.uk/projects/fastqc/
		reeBayes is a Bayesian genetic variant detector designed to find small polymorphisms, specifically SNPs (single-nucleotide polymorphisms), indets	
		(insertions and deletions), MNPs (multi-nucleotide polymorphisms), and complex events (composite insertion and substitution events) smaller than the	
123	freebayes	length of a short-read sequencing alignment.	http://bioinformatics.bc.edu/marthlab/wiki/index.php/FreeBayes
124	lammps	LAMMPS is a classical molecular dynamics code, and an acronym for Large-scale Atomic/Molecular Massively Parallel Simulator.	http://lammps.sandia.gov
		NAMD, recipient of a 2002 Gordon Bell Award and a 2012 Sidney Fernbach Award, is a parallel molecular dynamics code designed for high-performance	
		simulation of large biomolecular systems. Based on Charm++ parallel objects, NAMD scales to hundreds of cores for typical simulations and beyond	
125	namd	200,000 cores for the largest simulations.	http://www.ks.ujuc.edu/Research/namd/
126	raxml	RAVMI Version 8: A tool for Phylogenetic Analysis and Post-Analysis of Large Phylogenies	http://sco.h-its.org/exelixis/web/software/raxml/index.html
127	samtools	SAM (Sequence Alignment/Map) format is a generic format for storing large nucleotide sequence alignments	http://camtools.surreforme.net
120	schrodinger		http://darh.knime.org/forum/schroadinger.extensions/announcing.schroadinger.software.relasse 2014.2.0
120	somounger	The Sequence Read Archive (SPA) stores raw equence data from "part generation" equencing technologies including 454 tertarrent. Illuming SOLID	mp.//tori.kimic.org/forum/son/beuinger-extensions/announomy-son/beuinger-son/ware-release-2014-2-0
100	arataalkit	The organic mean many many lots is an sequence data from new yeneration sequencing technologies including 434, tornorrent, illumina, SOLID, Unices and Complete Life Sections	http://www.nahi.alm.aik.gov/Troaco/ara/ara/ara.aci/wiew.acifuara
129	STALUUIKIT	melicos ano complete Lile Sciences.	http://www.ncob.nim.nin.gov/iraces/sra/sra.cgi/view=sonware
130	topnat	TOPITAL IS A TASE SPIRE JUICTION MAPPER FOR KINA-SEQ FEADS.	nttp://cco.jnu.edu/softWare/tophat/index.shtml
1			
131	vmd	WMD is a molecular visualization program for displaying, animating, and analyzing large biomolecular systems using 3-D graphics and built-in scripting.	http://www.ks.uiuc.edu/Research/vmd/
132	voro++	Voro++ is a software library for carrying out three-dimensional computations of the Voronoi tessellation.	http://math.lbl.gov/voro++/
133	absorb	This software package implements ABSORB: Atlas Building by Self-Organized Registration and Bundling	http://www.nitrc.org/projects/absorb/
		ABySS is a de novo, parallel, paired-end sequence assembler that is designed for short reads. The single-processor version is useful for assembling	
134	abyss	genomes up to 100 Mbases in size. The parallel version is implemented using MPI and is capable of assembling larger genomes.	http://www.bcgsc.ca/platform/bioinfo/software/abyss
<u> </u>	1	AFNI (which might be an acronym for Analysis of Functional NeuroImages) is a set of C programs for processing analyzing and displaying functional MRI	
135	afni	EMBI data - a technique for manning human brain activity	http://afni.nimh.nih.gov/afni/
120	amos	AMOS is a collection of tools for anome assembly	http://amore.coursectore.o.pot
100	atros	ramedo is a concernent or consister generation and win from biological magnetization	http://www.ombl.bomburg.do/biogoove/coftware.btml
13/	aisas	n program some for sman-angle scattering data analysis nom biological macromolecules	nttp://www.enurnanbulg.ue/biosaxs/software.ntm
		AU LOJUEW IS an automated image reconstruction system that coordinates the execution of the parallel, numerically-intensive codes P3DR, PCUT, POR,	
138	auto3dem	PPFI, and PSF.	nttp://cryoem.ucsd.edu/programs.shtm
1		The Bcl2FastQ conversion software is a tool to handle bcl conversion and demultiplexing. Version 1.8.4 has added ability to mask multiple adapter	
1		sequences per read, has standard Illumina adapter sequences included in the bcl2fastq installation, and the stringency of the adapter masking feature is	
139	bcl2fastq	now configurable.	http://support.illumina.com/downloads/bcl2fastg_conversion_software_184.html
		BEDOPS is an open-source command-line toolkit that performs highly efficient and scalable Boolean and other set operations, statistical calculations.	
1		archiving, conversion and other management of genomic data of arbitrary scale. Tasks can be easily solit by chromosome for distributing whole-genome	
140	bedops	analyses across a computational cluster	http://bedops.readthedocs.org/en/latest/
140		Binawk is a antensity of the second	mp.// bodopol.oudulodos.org/orinates/
1.44	biopult	End where EACE A() and TAB relimined to any advergence support or several common biological data formats, including optionally gzip ed BED, GFF,	https://www.biostore.org/c/47754/
141	DIDAWK	SAW, VCF, FASTAVQ and FAS-deminied formats with course names.	nttps://www.biostans.org/p/4/1/51/
I		BLAT on DIVA is designed to quickly find sequences of 95% and greater similarity of length 25 bases or more. It may miss more divergent or shorter	
142	blat	sequence alignments. It will find perfect sequence matches of 20 bases.	http://genome.ucsc.edu/cgi-bin/hgBlat
1		Bsoft is a collection of programs and a platform for development of software for image and molecular processing in structural biology. Problems in	
1		structural biology are approached with a highly modular design, allowing fast development of new algorithms without the burden of issues such as file I/O.	
143	bsoft	It provides an easily accessible interface, a resource that can be and has been used in other packages.	http://lsbr.niams.nih.gov/bsoft/bsoft.html
144	cd-hit	CD-HIT is a very widely used program for clustering and comparing protein or nucleotide sequences.	http://weizhong-lab.ucsd.edu/cd-hit/
145	cnypator	a tool for CNV discovery and genotyping from denth of read mapping	http://sv.gersteinlab.org
170	onnator	a teorier erve accevery and generyping nom departer read mapping.	initial and the second s

	А	S	Т
-	<i></i>		
		cutadapt is used to remove adapter sequences from high-throughput sequencing data. This is usually necessary when the read length of the sequencing	
146	cutadapt	machine is longer than the molecule that is sequenced, for example when sequencing microRNAs.	https://sites.google.com/a/brown.edu/bioinformatics-in-biomed/cutadapt
147	dwgsim	Whole Genome Simulator for Next-Generation Sequencing	https://github.com/nh13/DWGSIM
		The EIGENSOFT package combines functionality from our population genetics methods (Patterson et al. 2006) and our EIGENSTRAT stratification	
148	eigensoft	correction method (Price et al. 2006)	http://genetics.med.harvard.edu/reich/Reich Lab/Software.html
		ESTScan is a program that can detect coding regions in DNA sequences, even if they are of low guality. ESTScan will also detect and correct sequencing	
149	estscan	errors that lead to frameshifts.	http://estscan.sourceforge.net
150	fastr	The EASTX-Toolkit is a collection of command line tools for Short-Reads EASTA/EASTO files preprocessing	http://bapponlab.csh.edu/fastx_toolkit/index.html
151	froorurfor	An onen source software suite for processing and analyzing (human) brain MRI images	http://fameurica.net
150	fel	An open source software softe for processing and analyzing (name) bran men magness.	http://necsus.dokion.net/n/co/fal_5_0.complete.html
132	151	FSL is a complemensive library or image analysis and statistical tools for highly, with and off brain imaging data.	http://neuro.debian.nevpkgs/nsi-5.0-complete.num
153	gmap-gsnap	GMAP: A Genomic Mapping and Alignment Program for mRNA and EST Sequences, and GSNAP: Genomic Short-read Nucleotide Alignment Program	http://research-pub.gene.com/gmap/
		GenomeTools genome analysis system is a set of bioinformatics tools (in the realm of genome informatics) combined into a single binary. more details at	
154	gt	GenomeTools.	https://wiki.gacrc.uga.edu/wiki/GenomeTools
		GView is a Java package used to display and navigate bacterial genomes. GView is useful for producing high-quality genome maps for use in publications	
155	gview	and websites, or as a visualization tool in a sequence annotation pipeline.	https://www.gview.ca/wiki/GView/WebHome
		HMMER is used for searching sequence databases for homologs of protein sequences, and for making protein sequence alignments. It implements	· · ·
156	hmmer	methods using probabilistic models called profile bidden Markov models (profile HMMs)	http://hmmer.janelia.org/software
		The Integrative Life Sciences Viewer (IGV) is a high-performance visualization tool for integrative exploration of large integrated genomic datasets. It	
157	iov	integration of the second second second and part apparently compared and apparently compared attained and apparently of apparently of a second and part apparently of apparently of apparently of a second and part apparently of apparently of apparently of a second apparently of apparently of apparently of a second apparently of a second apparently of apparently of apparently of a second ap	http://www.broadingtitute.org/au/
137	igv	supports a vide valiety of data types, including analybased and including relation sequence data, and genome amobations.	nip.//www.broadinsulate.org/igw
		The integrative Life Sciences viewer (IGV) is a high-performance visualization tool for interactive exploration of large, integrated genomic datasets. It	
158	Igvtoois	supports a wide variety or data types, including array-based and next-generation sequence data, and genomic annotations.	nttp://www.broadinstitute.org/igv/
159	Idnat	LUDIAL IS A PACKAGE WRITTEN IN THE C and C++ languages for the analysis of recombination rates from population genetic data.	http://dnat.sourcetorge.net
160	Idhelmet	LDhelmet is a sottware program for statistical inference of fine-scale crossover recombination rates from population genetic data.	http://sourcetorge.net/projects/ldhelmet/
161	meme	MEME is a tool for discovering motifs in a group of related DNA or protein sequences.	http://ebi.edu.au/ftp/software/MEME/
162	migrate	Migration and gene flow estimator	http://popgen.sc.fsu.edu/Migrate/Migrate-n.html
163	mothur	This project seeks to develop a single piece of open-source, expandable software to fill the bioinformatics needs of the microbial ecology community.	http://www.mothur.org
		Mugsy is a multiple whole genome aligner. Mugsy uses Nucmer for pairwise alignment, a custom graph based segmentation proceeding for identifying	
		collinear regions, and the segment-based progressive multiple alignment strateging from sages "Torfae Museu segment-based regiment in the form of multi-	
164		common regions, and the segment based progressive multiple angument strategy non sequit. Founde, mugsy accepts that genomes in the form of multi-	http://www.courseferee.pet
104	muysy	r Ao Ines and ques not require a reference genome.	nup.//mugsy.sourcerorge.net
		MUMmer is a system for rapidly aligning entire genomes, whether in complete or draft form. For example, MUMmer 3.0 can find all 20-basepair or longer	
165	mummer	exact matches between a pair of 5-megabase genomes in 13.7 seconds, using 78 MB of memory, on a 2.4 GHz Linux desktop computer.	http://mummer.sourceforge.net
166	pandaseq	PAired-eND Assembler for DNA sequences	https://github.com/neufeld/pandaseg
		PASA, acronym for Program to Assemble Spliced Alignments, is a eukaryotic genome annotation tool that exploits spliced alignments of expressed	
		transcript sequences to automatically model gene structures, and to maintain gene structure annotation consistent with the most recently available	
167	pasa	experimental sequence data.	http://pasa.sourceforge.net
.07	pada	RepeatMasker is a program that screens DNA sequences for interspersed repeats and low complexity DNA sequences. The output of the program is a	mp.//padicouroologo.net
		detailed appendix to a program that below to recent in the appendix on probability of the appendix of the appendix to the appendix of the appe	
160	repeatmosker	detailed antidation of the repeats that are present in the query sequence as well as a modified version of the query sequence in which all the antidated	http://www.represtreader.org
168	repeatmasker	repeats have been masked (derault: replaced by Ns).	nttp://www.repeatmasker.org
169	rtax	Taxonomic classifications of single- and pared-end sequences with RTAX	http://qime.org/tutorials/rtax.html
170	sailtish	Rapid Alignment-free Quantification of Isoform Abundance	http://www.cs.cmu.edu/~ckingst/software/sailtish/README.html
		Seqtk is a fast and lightweight tool for processing sequences in the FASTA or FASTQ format. It seamlessly parses both FASTA and FASTQ files which	
171	seqtk	can also be optionally compressed by gzip.	https://wiki.gacrc.uga.edu/wiki/Seqtk
		SHRiMP is a software package for aligning genomic reads against a target genome. It was primarily developed with the multitudinous short reads of next	
172	shrimp	generation sequencing machines in mind, as well as Applied Biosystem's colourspace genomic representation.	http://compbio.cs.toronto.edu/shrimp/
173	smalt	SMALT efficiently aligns DNA sequencing reads with a reference genome.	https://www.sanger.ac.uk/resources/software/smalt/
174	smrtanalysis	PacBio's open source software suite for single molecule, real-time sequencing	http://www.pach.com/devnet/
	omitanaiyolo	SAKE is a da novo assemblar for short DNA sequence reads. It is designed to belo leverage the information from short sequences reads by assembling	mp.//www.pado.com/domov
175	oooko	Some is a service and experience to short prive sequence reads, in a designed to help reverge the information from short sequences reads by assembling	http://www.bases.cs/platfarm/bicipfs/asftware/asska
175	SSake	mem mito condes and scanolos mar can be used to characterize novel sequencing targets.	http://www.bcgsc.ca/platform/bioinfo/software/ssake
		Stacks is a software pipeline for building four from short-read sequences, such as those generated on the information. Stacks was developed to work	
		with restriction enzyme-based data, such as RAD-seq, for the purpose of building genetic maps and conducting population Life Sciences and	
176	STACKS	pnyiogeography.	http://creskolab.uoregon.edu/stacks/
177	structure	Ine program structure is a tree software package tor using multi-locus genotype data to investigate population structure.	http://pritchardlab.stanford.edu/structure.html
		SAM (Sequence Alignment/Map) is a flexible generic format for storing nucleotide sequence alignment. SAM tools provide efficient utilities on	
178	tabix	manipulating alignments in the SAM format.	http://samtools.sourceforge.net
179	trimmomatic	Trimmomatic: A flexible read trimming tool for Illumina NGS data	http://www.usadellab.org/cms/index.php?page=trimmomatic
180	velvet	Velvet is a de novo genomic assembler specially designed for short read sequencing technologies	https://www.ebi.ac.uk/~zerbino/velvet/
		Xmipp is a specialized suite of image processing programs, primarily aimed at obtaining the 3D reconstruction of biological specimens from large sets of	
181	adimx	projection images acquired by transmission electron microscopy	http://xmipp.cnb.csic.es/twiki/bin/view/Xmipp/WebHome
100	abaque	Abarus EEA. (formerly ABAOLIS) is a software suite for finite element analysis and computer-aided Physical Sciences, originally released in 1079	http://www.cimulia.com
102	avayas	ANSYS for provide roll word uplus to the Dhuring Science dorign present any and the analysis and computer and a sound of the the Dhuring Sciences. Design present and any and the to the Dhuring Science dorign present and an any and the analysis and computer and any and the science of the science.	Intpartwww.sinuid.colli
400		rave to technology acts rearmond value to the Englished Sciences design process. Robust simulation capabilities codpied with a flexible, adaptive	
183	ansys	and meteouse deriver emodelity and univerning value.	http://www.ansys.com
184	campañ	Molecular Modeling SoftWare	nttp://campan.sourcetorge.net
185	comsol	Simulation Tool for Electrical, Mechanical, Fluid Flow, and Chemical applications	http://www.comsoi.com
186	dl_poly	DL_POLY is a general purpose classical molecular dynamics (MD) simulation software developed at Daresbury Laboratory by I.T. Todorov and W. Smith.	http://www.sttc.ac.uk/SCD/research/app/ccg/software/DL_POLY/44516.aspx
187	dl_poly_classic	DL_POLY Classic is a general purpose molecular dynamics simulation package developed at Daresbury Laboratory	http://www.ccp5.ac.uk/DL_POLY_CLASSIC/
188	Extend-Project	Community contributed extensions to the OpenFOAM CFD toolbox, not approved or endorsed by ESI Group	http://www.extend-project.de/
189	gaussian	Gaussian 09 is the latest in the Gaussian series of programs. It provides state-of-the-art capabilities for electronic structure modeling.	http://www.gaussian.com
	-	LS-DYNA is a general-purpose finite element program capable of simulating complex real world problems. It is used by the automobile, aerospace.	
190	Isdvna	construction, military, manufacturing, and bioPhysical Sciences industries.	http://www.lstc.com/products/ls-dvna
		Materials Studio ® is a software environment that brings the world's most advanced and validated materials simulation technology to desktop computing	
101	Materials Studio	solving key problems throughout the R&D process.	https://wiki.csiro.au/display/ASC/Accelrys+Materials+Studio
131			mpos/ millionio.co/orophy/ noo/noon/coloring millionas rotadio
100	nucham	APMCI MDI is a completely rewritten implementation of the APMCI and sided communication interface that uses MDI DMA for one olded communication	http://wiki.mpick.org/organi.http://www.puphorg.ou.org/index.php/Douplood
192	ananfaam	non-mining a complexity rewritten implementation of the artivity of ensured communication interface that uses with twink for one-sided communication.	http://wikianpior.org/afficientp://www.nwcnem-sw.org/index.pnp/Download
193	openidam	Openic Own is a nee, open source CFD sortware package developed by OpenCFD Ltd at ESI Group and distributed by the OpenFOAM Foundation	nup.//www.openroam.org
		Pointwise is our leading mesh generation software product for CFD, encompassing nearly three decades of technology development. The product	
194	pointwise	Iteratives a modern software framework and workflow and provides tools like 64-bit support, multi-threading, and Windows/Linux/Mac portability.	http://www.pointwise.com
		Q-Chem is a comprehensive ab initio quantum Physical Sciences package for accurate predictions of molecular structures, reactivities, and vibrational,	
195	qchem	electronic and NMR spectra.	http://www.q-chem.com
		Quantum Expresso (QE) is an integrated suite of Open-Source computer codes for electronic-structure calculations and materials modeling at the	
196	qe	nanoscale. It is based on density-functional theory, plane waves, and pseudopotentials.	http://www.quantum-espresso.org/download/
197	quantumwise	Simulation software for panoscience professionals	http://quantumwise.com

	Δ	2	T
100			
198	starccm+	StarcCM+ is entire Physical Sciences process for solving problems involving flow (of fluids or solids), neat transfer and stress.	nttp://www.cd-adapco.com/products/star-ccm®
		Tecplot software empowers engineers and scientists working with Physical Sciences (CFD) to discover, analyze, and understand information in complex	
199	tecplot	data.	http://www.tecplot.com/
		The Trilinos Project is an effort to develop algorithms and enabling technologies within an object-oriented software framework for the solution of large-	
200	trilinos	scale, complex multi-Physical Sciences Physical Sciences and scientific problems.	http://trilings.org
200	dimited	The Vienne Ah initia Simulation Deckage (VASD) is a computer program for atomic scale materials modelling and electronic structure calculations and	independient of the second s
		The vienna Ab initio Simulation Package (VASP) is a computer program for atomic scale materials modelling, e.g. electronic structure calculations and	
201	vasp5	quantum-mechanical molecular dynamics, from first principles.	nttp://www.vasp.at
202	xcrysden	Construction and visuazliation of atomistic structures and properties	http://www.xcrysden.org
		ABINIT is a package whose main program allows one to find the total energy, charge density and electronic structure of systems made of electrons and	
203	abinit	nuclei (molecules and periodic solids) within Density Functional Theory (DET), using oseudopotentials and a planewave or wavelet basis	http://www.abipit.org
200	abilit	TAT is a apprised to find to find the collection of allow hears tools (0117), alloy pooled option and a planoward of wardoor backs.	http://www.abincorg
204	alal	ATAT is a genetic name mathematic a concernor of any free yours	http://audm.mit.edu/www/audw/atat
205	cantera	Cantera is a suite or object-oriented software tools for problems involving chemical kinetics, thermodynamics, and/or transport processes.	http://www.cantera.org/docs/spninx/ntmi/index.ntmi
206	eman2		http://ncmi.bcm.edu/ncmi/software/counter_222/software_122
		GROMACS is a versatile package to perform molecular dynamics, i.e. simulate the Newtonian equations of motion for systems with hundreds to millions of	
207	gromacs	particles.	http://www.gromacs.org
		Meep (or MEEP) is a free finite-difference time-domain (FDTD) simulation software package developed at MIT to model electromagnetic systems, along	
208	meen	with our MPB eigenmode package. I	http://ab-initio.mit.edu/wiki/index.php/Meep
200	moop	Maldon is a package for disclosing Malagular Density from the Ab Initia packages CAMESS LIK CAMESS LIS and CAUSSIAN and the Sami Empirical	n na maamaada maamada phi moop
		wolden is a package for displaying wolecular bensity form the Ab Initio packages GAMIEGSON, GAMIEGSOS and GAOSGIAN and the Semi-Empirical	
209	molden	packages Mopac/Ampac, it also supports a number of other programs via the Molden Format.	http://www.cmbi.ru.nl/molden/
		NASTRAN is a finite element analysis (FEA) program that was originally developed for NASA in the late 1960s under United States government funding	
210	nastran	for the Aerospace industry.	http://www.mscsoftware.com/product/msc-nastran, http://en.wikipedia.org/wiki/Nastran
		Nmag is a micromagnetic simulation package. It has been developed at the University of Southampton with substantial contributions from Hans Fangohr,	
211	nmag	Thomas Fischbacher, Matteo Franchin, It is released under the GNU GPL	http://nmag.soton.ac.uk/nmag/0.2/index.html
		Onen Babel is a chemical tophox designed to speak the many languages of chemical data. It's an onen, collaborative project allowing anyone to search	
240	ananhahal	open sate in a similar tonice designed to speak the many rangedes of definition data. It's an open, contact end for the set of the form realized and the set of the s	http://openhabel.org/wiki/Main_Dogo
212	openbabel	convert, analyze, or store data from molecular modeling, Physical Sciences, solid-state materials, bioPhysical Sciences, or related areas.	nttp://openbabel.org/Wiki/Main_Page
		Patran is the world's most widely used pre/post-processing software for Finite Element Analysis (FEA), providing solid modeling, meshing, analysis setup	
213	patran	and post-processing for multiple solvers including MSC Nastran, Marc, Abaqus, LS-DYNA, ANSYS, and Pam-Crash.	http://www.mscsoftware.com/product/patran
214	phenix	PHENIX is a software suite for the automated determination of molecular structures using X-ray crystallography and other methods.	http://www.phenix-online.org
		PHOEBE stands for Physical Sciences Of Eclipsing Binaries. It is a tool for modeling eclipsing binary stars based on photometric, spectroscopic	
215	phoebe	interferometric and/or polarimetric data	http://phoebe-project.org
213	p.10000	monorman and postminute data.	mps/procedo projekterg
		RELICIN (for REgularised Likelinood Optimisation, pronounce rely-on) is a stand-alone computer program that employs an empirical Bayesian approach to	
216	relion	refinement of (multiple) 3D reconstructions or 2D class averages in electron cryo-microscopy (cryo-EM).	http://www2.mrc-lmb.cam.ac.uk/relion/index.php/Main_Page
217	unitah	The Uintah software suite is a set of libraries and applications for simulating and analyzing complex chemical and physical reactions.	http://www.sci.utah.edu/software/uintah.html
		WanT is a GPL open-source scientific software aimed at simulating quantum transport properties (conductance) of nanojunctions (atomistically described)	
218	WanT	by using a real space Wannier function description of the Hamiltonian of the system.	http://www.wannier-transport.org/wiki/index.php/Main_Page
		Vamba is a FORTRAN/C code for Many-Body calculations in solid state and molecular Physical Sciences. Vamba can calculate in an Ab-Initio manner	
210	vombo	Tambo is a contribute of a constraint of the solution in solutions in solution and the contribute of the activity of the solution in the solution of the solut	http://www.wombo.co.do.org
219	yambo	absolption, energy loss and quasipancies properties.	ntp://www.yambo-code.org
220	R (+all the modules)	The R Project Statistical Computing Package	http://www.r-project.org
		The AMPL system is a sophisticated modeling tool that supports the entire optimization modeling lifecycle: development, testing, deployment, and	
221	ampl	maintenance.	http://ampl.com
222	cplex	High-performance mathematical programming solver for linear programming, mixed integer programming, and guadratic programming	http://www-01.ibm.com/software/commerce/optimization/cplex-optimizer/
		The General Algebraic Modeling System (GAMS) is a high-level modeling system for mathematical programming and optimization. It consists of a	
		language compiler and a stable of integrated high-performance solvers. GAMS is tailored for complex large scale modeling applications, and allows you	
222	20m2	tangado comprior are o anto a magnato ingri portornario do torio o anto to ano to comprov, argo codo modorny approximento, and anoro you	http://www.como.com
223	yanis	to build raige maintainable models that can be adapted quickly to new situations.	nto.//www.gans.com
224	gauss		http://www.aptech.com
225	gurobi	Gurobi Optimizer 5.6 - State-of-the-Art Mathematical Programming Solver	nttp://www.gurobi.com
		KNITRO is an ideal tool for solving problems, such as pricing or revenue management, parameter identification in financial and industrial models or design	
226	knitro	and operation of electricity and gas transportation networks.	http://www.artelys.com/en/optimization-tools/knitro
		Mathematica is a computational software program used in many scientific, Physical Sciences, mathematical and computing fields, based on symbolic	
227	mathematica	mathematics	http://www.wolfram.com/mathematica/
220	matlah	MATLADE is a bigh level language and interactive environment for numerical computation, visualization, and pregramming	nue motivere com
220	manad	men carbon is a nymetrix i anguage and interactive environment on numerical Computation, Visualization, and programming.	www.induiwoirs.com
		raraview is an open-source, multi-platform data analysis and visualization application. Paraview users can quickly build visualizations to analyze their	
229	paraview	data using qualitative and quantitative techniques.	nttp://www.paraview.org/
230	sas		http://www.sas.com/
		Stata statistical software is a complete, integrated statistical software package that provides everything you need for data analysis, data management, and	
231	stata	graphics.	http://www.stata.com
222		The TOMLAB Optimization Environment is a powerful optimization platform and modeling language for solving applied optimization problems in MATLAP	http://tomont.com/tomlab/
202	0.470	Aviza is a Dispanization Environment is a power or optimization planet and modeling language for solving applied optimization problems in MATEAD.	http://unicpit.com/onfluors/puizo2d/
200	anto	ranzo is a operatorise sontwate for soletimite and industrial data.	Intp://www.noi.com/collwalC/avi2000/
		The graphical models toolkit (GMTK) is an open source, publicly available toolkit for rapidly prototyping statistical models using dynamic graphical models	
234	gmtk	(DGMS) and dynamic Bayesian networks (DBNs).	nttp://meiodi.ee.washington.edu/gmtk/
		Macaulay2 is a software system devoted to supporting research in algebraic geometry and commutative algebra, whose creation has been funded by the	
235	macaulay2	National Science Foundation since 1992.	http://www.math.uiuc.edu/Macaulay2/
236	maple	Maple combines the world's most powerful mathematical computation engine with an intuitive, clickable user interface.	http://www.maplesoft.com/products/maple/
237	openbugs	BUGS is a software package for performing Bayesian inference Using Gibbs Sampling	http://www.openbugs.net/w/FrontPage
239	oprofile	OProfile is an open source protect that includes a statistical profiler for Linux systems, consolie of profiling all running code at low everteed	http://oprofile.sourceforce.pet/about/
200	opromo	The DDB closeling is a point of the provide a statistical profile for the track systems, depaids of profiling an furning doub at the operational statistical profile the systems and a force of the system of the sy	mp//opromotocial.gomerabodu
000	ada alaasifaa	The row Gassine is a name payesian classiner that can rapidly and accurately provides taxonomic assignments from domain to genus, with confidence	
239	rap_classifer	esumates for each assignment.	nttps://rap.cme.msu.eau
		Sage is a tree open-source mathematics software system licensed under the GPL. It builds on top of many existing open-source packages: NumPy, SciPy,	
240	sage	matplotlib, Sympy, Maxima, GAP, FLINT, R and many more.	http://www.sagemath.org
		Slic3r is the tool you need to convert a digital 3D model into printing instructions for your 3D printer. It cuts the model into horizontal slices (lavers).	
2/1	slic3r	generates toolpaths to fill them and calculates the amount of material to be extruded	http://slic3r.org
241	0.000	IDM SDES Statistics is an interacted family of products that addresses the antiro analytical process from alapping to date collection to execution	mps/sitering
o		now or so statistics is an integrated ramity or products that addresses the entire analytical process, from planning to data collection to analysis, reporting	
242	spss	and deproyment.	nttp://www-u1.ibm.com/software/analytics/spss/products/statistics/
243	INTERACTIVE - HAMMER		
		The Portable Hardware Locality (hwloc) software package provides a portable abstraction of the hierarchical topology, including NUMA memory nodes,	
		shared caches, processor sockets, processor cores and processing units (logical processors or "threads"). It primarily aims at helping applications with	
244	hwloc	gathering information about modern computing hardware so as to exploit it accordingly and efficiently.	http://www.open-mpi.org/projects/hwloc/
245	mumps	MUMPS: a MUltifrontal Massively Parallel sparse direct Solver	http://mumps.enseeiht.fr/
240	at	is a proper platform policition and III from our for dougleport using CLL or CML a CCC % laugCariet like language	http://ntmpsciedcontcli/
240	yı metrodi e	is a cross-pration application and on trainework for developers using C++ or QML, a CSS & JavaScript like language	Inter//recipioject.org/
247	Istuuid	Rotucio IDE is a poweriul and productive user interface for R	nttp://www.isituaio.com/
	aloog ppl	(1) oof is a software which generates loops for scanning 7-nolyhedra	https://packages.debiap.org/sid/devel/cloog-ppl

	A	S	Т
		ANSYS Fluent software contains the broad physical modeling capabilities needed to model flow, turbulence, heat transfer, and reactions for industrial	
		applications ranging from air flow over an aircraft wing to combustion in a furnace, from bubble columns to oil platforms, from blood flow to semiconductor	
249	fluent	manufacturing, and from clean room design to wastewater treatment plants	http://www.apsys.com/Products/Simulation+Technology/Eluid+Dynamics/Eluid+Dynamics+Products/ANSYS+Eluent
250	aridaen	Gridgen is Pointwise's lengery software used by engineers and scientists worldwide since 1984 to reliably generate high guality grids for CED	http://www.pointwise.com/arideen/
251	hfss	ANSYS HESS software is the industry standard for simulation 3-D full-wave electromagnetic fields	http://www.apsys.com/Products/Simulation+Technology/Electronics/Signal+Integrity/ANSYS+HESS
		The ANSYS Maxwell course will teach the students how to design and analyze 2D and 3D electromagnetic and electromechanical devices such as	
252	maxwell	motors actuators transformers sensors and coils	http://www.ansvs.com/Support/Training+Center/Courses/Introduction+to+&NSVS+Maxwell+2D-3D
253	siwave	ANSYS Stearers, a specialized design platform for power integrity, signal integrity and EMI analysis of electronic packages and PCBs	http://www.ansys.com/Products/Simulation+Technology/Electronics/Signal+Intertity/ANSYS+Slwave
200	omaro	The SPM software because a been designed for the analysis of brain imaging data sequences. The sequences can be a series of images from different	maps/www.tanejo.com/reddold/ormalation/recombining/jichodromod/orgina/rintegrity/wite/re-re-ormate
254	som	cohorts, or time-series from the same subject. The current release is designed for the analysis of fMRI. PET. SPECT. EEG and MEG.	http://www.fil.ion.ucl.ac.uk/spm/
		Adams is the most widely used multibody dynamics and motion analysis software in the world. Adams helps engineers to study the dynamics of moving	
255	Adams	parts, how loads and forces are distributed throughout mechanical systems, and to improve and optimize the performance of their products.	http://www.mscsoftware.com/adams-student-edition
		DINA System for linear and ponlinear finite element analysis of solids and structures, heat transfer. CED and electromagnetics. ADINA also offers a	
256	Adina	comprehensive array of multiphysics capabilities including fluid-structure interaction and thermo-mechanical coupling	http://www.adina.com/index.shtml
		ANTs extracts information from complex datasets that include imaging (Word Cloud). Paired with ANTsR (answer). ANTs is useful for managing.	
		interpreting and visualizing multidimensional data. ANTs is popularly considered a state-of-the-art medical image registration and segmentation toolkit.	
		ANTsR is an emerging tool supporting standardized multimodality image analysis. ANTs depends on the Insight ToolKit (ITK), a widely used medical	
257	Advanced Normalized Tools (ANTs)	image processing library to which ANTs developers contribute.	http://stnava.github.io/ANTs/
258	Ansoft Designer	2	http://www.ansvs.com/Support/Platform+Support/ch.Ansoft+Products+14.0.hu
259	Art		
		The Atomistic Simulation Environment (ASE) is a set of tools and Python modules for setting up, manipulating, running, visualizing and analyzing atomistic	
260	Atomic Simulation Environment (ASE)	simulations. The code is freely available under the GNU LGPL license.	https://wiki.fysik.dtu.dk/ase/
261	BCT		
262	Beagle		
		BEAST is a cross-platform program for Bayesian analysis of molecular sequences using MCMC. It is entirely orientated towards rooted, time-measured	
		phylogenies inferred using strict or relaxed molecular clock models. It can be used as a method of reconstructing phylogenies but is also a framework for	
		testing evolutionary hypotheses without conditioning on a single tree topology. BEAST uses MCMC to average over tree space, so that each tree is	
		weighted proportional to its posterior probability. We include a simple to use user-interface program for setting up standard analyses and a suit of	
263	Beast	programs for analysing the results.	http://beast.bio.ed.ac.uk/
264	Blender		
1		CCP4 exists to produce and support a world-leading, integrated suite of programs that allows researchers to determine macromolecular structures by X-	
		ray crystallography, and other biophysical techniques. CCP4 aims to develop and support the development of cutting edge approaches to experimental	
265	ccp4	determination and analysis of protein structure, and integrate these approaches into the suit	http://www.ccp4.ac.uk/about.php
266	charmm	is a versatile and widely used molecular simulation program with broad application to many-particle systems	http://www.charmm.org/
		UCSF Chimera is a highly extensible program for interactive visualization and analysis of molecular structures and related data, including density maps,	
267	chimera	supramolecular assemblies, sequence alignments, docking results, trajectories, and conformational ensembles	https://www.cgl.ucsf.edu/chimera/index.html
268	chipmonk	ChIPMonk is a program to enable the visualisation and analysis of ChIP-on-chip data	http://www.bioinformatics.babraham.ac.uk/projects/chipmonk/
		Clustal Omega is the latest addition to the Clustal family. It offers a significant increase in scalability over previous versions, allowing hundreds of	
	1	thousands of sequences to be aligned in only a few hours. It will also make use of multiple processors, where present. In addition, the quality of	
269	ciustal	alignments is superior to previous versions, as measured by a range of popular benchmarks.	nttp://www.clustai.org/omega/
		Crystallography & NMR System (CNS) is the result of an international collaborative effort among several research groups. The program has been	
		designed to provide a next the multi-level interactinical approach for the most commonly used algorithms in macromolecular structure determination.	
270	ana askus	Highlights include neavy atom searching, experimental phasing (including MAD and MiR), density modification, crystallographic refinement with maximum likelihead torgate, and NMD extensive sequences a locuring block density modification equilated atom	
270	cris_solve	intelinood targets, and wink structure calculation using NOEs, 5-coupling, chemical shint, and oppilar coupling data	ntp.//cns-online.org/
271	comet	CompEcon is a set of MATLAP functions for solving a variety of problems in accomptise and finance. The library functions include radifieding and	
		Completion is a set of market distributions for solving a variety of problems in economics and initiative. The initiative functional formation and and a set of market distribution is not dependent of an economic and and market distribution and other functional formation and and a set of market distribution and and a set of market distribution and and and a set of market distribution and and and and a set of market distribution and and and and and and and and and an	
		opunitzation routines for anergial functions and for common probability distributions, general solutions for denserial functions and for common probability distributions and are common probability distributions are common probability distributions and are common probability distributions are common probability distributing and are common probability distributions	
		boundary value problems) multiples for solving discrete and continuous time dynamic programming problems, and a general solver for financial derivatives	
272	compecon	boundary value proteins), realities of solving discrete and continuous time dynamic programming proteins, and a general solver for internolated values of (honds futures onlines)	http://www4.ncsu.edu/~pfackler/compecon/toolbox.html
273	csim		
274	Data Analysis		
<u> </u>		Desmond is a software package developed at D. E. Shaw Research to perform high-speed molecular dynamics simulations of biological systems on	
		conventional commodity clusters. The code uses novel parallel algorithms and numerical techniques to achieve high performance and accuracy on	
275	desmond	platforms containing a large number of processors, but may also be executed on a single computer.	http://www.mybiosoftware.com/3d-molecular-model/1065
276	dmatlab	Distributed Matlab	
		EEGLAB is an interactive Matlab toolbox for processing continuous and event-related EEG, MEG and other electrophysiological data incorporating	
		independent component analysis (ICA), time/frequency analysis, artifact rejection, event-related statistics, and several useful modes of visualization of the	
277	eeglab	averaged and single-trial data.	http://sccn.ucsd.edu/eeglab/
278	em2em	Program to convert images (2-D images and 3-D volumes) from/to formats typically used in 3-D electron microscopy (3DEM).	https://www.imagescience.de/em2em.html
279	eman	EMAN is a suite of scientific image processing tools aimed primarily at the transmission electron microscopy community	http://blake.bcm.edu/emanwiki/EMAN1
		EVIDENDS is a ree Open Source software analysis package specially developed for the needs of the molecular biology (e.g. EMBnet) user community. The	
0000		sortware automatically copes with data in a variety of formats and even allows transparent retrieval of sequence data from the web. Also, as extensive	
280	emboss faatBHASE	inoranes are provided with the package, it is a platform to allow other scientists to develop and release software in true open source spirit.	nttp://emposs.sourcerorge.nevwhat/
281	IdSIFTAGE	The program rate managements methods for estimating naplotypes and missing genotypes from population SNP genotype data.	http://www.ce.bedie.com/express_idense_technologies/rastphase
282	teappv	EAPPy is a general purpose tinite element analysis program which is designed for research and educational use	http://www.ce.berkeley.edu/projects/reap/reappv/
202	ferret	i errer is an interactive computer installization and analysis environment designed to meet the needs of oceanographies and meleorologists analyzing large and complex origidad data sets it ruis on most libry systems, and on Windows VDMTW using X windows for disclose	http://ferret.pmel.poga.gov/Ferret/home
203	lenet	large and complex gnoded data sets. In this of midst only systems, and on windows AP/N179X dsing A windows for display.	ntp.//ener.pmei.noaa.gov/rene/nome
		vide to be to be to be a set of generating seamed, so many some may all the approximate the program which recognizes reach a patients in text, defines the function	
284	flex	yriady, into the require expressions for each rule. Whenever it finds a match it executes the corresponding C code	http://flex.sourceforge.pet/
204	10%	a cross-platform C++ GIII tookit for UNIXM. Interest names the Windows® and MacOS®X FI IX provides modern GII functionality without the	n pan becomo regenere
285	fitk	bloat and supports 3D graphics via OpenGL® and its built-in GLUT emulation.	http://www.fitk.org/index.php
286	frappe		
287	frealign	Frealign is a program for high-resolution refinement of 3D reconstructions from crvo-EM images of single particles	http://grigoriefflab.janelia.org/frealign
		FTGL is a free cross-platform Open Source C++ library that uses Freetype2 to simplify rendering fonts in OpenGL applications. FTGL supports bitmaps.	
288	ftgl	pixmaps, texture maps, outlines, polygon mesh, and extruded polygon rendering modes.	http://sourceforge.net/projects/ftgl/
289	ga		
290	gambit		
291	gc	?	http://www.hboehm.info/gc/gc_source/
		The GLPK (GNU Linear Programming Kit) package is intended for solving large-scale linear programming (LP), mixed integer programming (MIP), and	
292	alpk	other related problems. It is a set of routines written in ANSI C and organized in the form of a callable library.	https://www.gnu.org/software/glpk/

	А	S	Т
		GNAT is an Ada (including Ada 2005) toolsat, integrated into the GCC compiler system GNAT includes GPS, a visual IDE, a comprehensive toolsuite	
202	anot	Charlis an Ada (including Ada 2005) forset, integrated includes compiler system charliner dates of 0, a visual DE, a compiler historic toosaite	http://www.gov.org/ooftworg/goot/
293	gilai	Including a visual debugger, and a set of inbranes and bindings	http://www.ght.org/software/ghat/
		GPAW is a density-functional theory (DFT) Python code based on the projector-augmented wave (PAW) method and the atomic simulation environment	
294	gpaw	(ASE). It uses real-space uniform grids and multigrid methods, atom-centered basis-functions or plane-waves	https://wiki.tysik.dtu.dk/gpaw/
		Provides a robust and efficient collection of tools and libraries which support reading, writing, and manipulating an image in over 88 major formats	
295	Graphics Magick	including important formats like DPX, GIF, JPEG, JPEG-2000, PNG, PDF, PNM, and TIF	http://www.graphicsmagick.org/
		GRASS GIS, commonly referred to as GRASS (Geographic Resources Analysis Support System), is a free and open source Geographic Information	
		System (GIS) software suite used for geospatial data management and analysis, image processing, graphics and maps production, spatial modeling, and	
206	drass	vieualization	http://grass.osgeo.org/
230	giuss	VISCANZAROT	mp.//grass.osgco.org/
		Haskeli is an advanced purely-functional programming language. An open-source product or more than twenty years of cutting-edge research, it allows	
		rapid development of robust, concise, correct software. With strong support for integration with other languages, built-inconcurrency and parallelism,	
297	haskell	debuggers, profilers, rich libraries and an active community, Haskell makes it easier to produce flexible, maintainable, high-quality software.	https://www.haskell.org/haskellwiki/Haskell
298	hkl-2000	provides high quality software and hardware solutions for X-ray crystallography laboratories and synchrotron beam lines.	http://www.hkl-xray.com/download-instructions-hkl-2000
299	icewm	IceWM is a window manager for the X Window System (freedesktop, XFree86). The goal of IceWM is speed, simplicity, and not getting in the user's way.	http://www.icewm.org/
300	idba	is a practical iterative De Bruijn Graph De Novo Assembler for sequence assembly in bioinfomatics	http://www.mybiosoftware.com/assembly-tools/7062
		IMAGIC is a high end environment for the analysis of images, spectra and other multi-dimensional data-sets, IMAGIC's software package is aimed at	
301	imagic	processing (huge) data sets from (crvo-) electron microscopy, especially in the field of single particle analyses in Structural Biology,	https://www.imagescience.de/imagic.html
302	impute	IMPLITE is a program for estimating ("imputing") unobserved genotynes in SNP association studies	http://mathgen.stats.ox.ac.uk/impute/impute.html
303	instruct	The of 2 is a program for countaining (implaining) anobcorred gonorpoor in oral accordance address	mp///mangonodolo/adolo/impato/impato/imp
204	intlab	INITI AR is the Matleh tealbay for reliable computing and self validating algorithms	http://www.ti2.tu.borburg.do/intloh/
304	inticio	int ca sine matter toron for remarke computing and seri-variadaring appointing	http://www.do.te-haloorg.do/httdb/
205	iar	reprove a similaritor sorware to graphicary design and control rarge-scale reactified models. Simulations in reproduct control real-world devices in fear-time. Ique as the strended by reprovement and expense times and automatic interference to increase the sector strended by the strended by the sector strended by the se	http://opurpefores.pot/orsionte/inv/
305	lqr	can be extended by new neuron, and synapse types, and custom interfaces to hardware.	nttp://sourcerorge.nevprojects/df/
306	ITIK	Image Registration Toolkit	nttp://www.doc.ic.ac.uk/~df/SoftWare/
307	itksnap	ITK-SNAP is a software application used to segment structures in 3D medical images	http://www.itksnap.org/pmwiki/pmwiki.php
308	j3d	Java 3D API	https://java3d.java.net/binary-builds.html
309	imol	an open-source Java viewer for chemical structures in 3D	http://jmol.sourceforge.net/
310	jweb		
		LaTeX is a high-quality typesetting system; it includes features designed for the production of technical and scientific documentation. LaTeX is the de facto	
311	latex	standard for the communication and publication of scientific documents.	http://latex-project.org/ftp.html
312	lev00	user friendly packages for DFT codes	http://www.mth.kcl.ac.uk/~lev/codes/lev00/index.html
		The LONE pipeline is a free workflow application primarily aimed at computational scientists. With the LONE Pipeline, users can quickly create workflows	
313	loni/pipeline	that take advantage of all the gradeet tools available in peuroimaging genomics bioinformatics at	http://pipeline.http://cipeline.ht
214	loni/pipeline	martake advantage of an me greatest tools available in neuroimaging, genomes, bioimormates, etc	http://pipemics.ong/.ucs.edu/
314	IOTI/Server_library	MachQue is a programming likeau for the implementation of simulation table in machanics. Its source code is maintunvitten is Quu with parts is Duthon	http://pipeline.ioni.usc.edu/ean/server-duide/conniduration/
045		and Co. (under development)	
313	meensys	and Go (under development).	http://mechsys.nongine.org/mode.sintmi
217	metouoluot	Merculy oners a comprehensive range of tools for SD structure instantization and the exploration of crystal packing.	http://www.ccdc.cam.ac.uk/solutions/riceSoftware/rages/riceWercuty.aspx
317	metavervet	An extension of verver assembler to be hove metagenome assembly non-short sequence reads	http://metavervet.dna.bio.keio.ac.jp/
040		The metaward binner is a graphical binning tool that makes use of multivariate statistics of tetranucleotide frequencies and interpolated markov models. It	
318	metawatt	also performs taxonomic assessment or binning quality (via BLAST).	nttp://sourcerorge.net/projects/metawait/
319	mni		http://www.mccauslandcenter.sc.edu/mricro/
		MRIcro allows Windows and Linux computers view medical images. It is a standalone program, but includes tools to complement SPM (software that	
320	mricro	allows neuroimagers to analyse MRI, tMRI and PET images). MRIcro allows efficient viewing and exporting of brain images.	http://www.mccauslandcenter.sc.edu/mricro/mricro.html
		MRIcron is a user-friendly and efficient piece of software which aims to assist you in viewing medical imagery with multiple layering, supporting HDR, NII	
321	mricron	and VOI formats, thus enabling you to analyze scans, X-rays and other similar files with minimal effort.	http://www.mccauslandcenter.sc.edu/mricro/mricron/install.html
		MSC Nastran is the world's most widely used Finite Element Analysis (FEA) solver. When it comes to simulating stress, dynamics, or vibration of real-	
		world, complex systems, MSC Nastran is still the best and most trusted software in the world – period. Today, manufacturers of everything from parts to	
322	nastran	complex assemblies are choosing the FEA solver that is reliable and accurate enough to be certified by the FAA and other regulatory agencies.	http://www.mscsoftware.com/msc-nastran-student-edition
		NetLogo is a programmable modeling environment for simulating natural and social phenomena.NetLogo is particularly well suited for modeling complex	
323	netlogo	systems developing over time	http://ccl.northwestern.edu/netlogo/
324	nltk	Natural Language Toolkit (NLTK) is a Python package for processing natural language text.	http://www.nltk.org/install.html
		nMOLDYN is an interactive analysis program for Molecular Dynamics simulations. It is especially designed for the computation and decomposition of	
325	nmoldyn	neutron scattering spectra, but also computes other quantities.	http://dirac.cnrs-orleans.fr/plone/software/nmoldvn/
-		numactl is an open source command-line software that implements support for simple NUMA policy. The NUMA (Non-Uniform Memory Access) policy	· · · · · · · · · · · · · · · · · · ·
326	numacti	allows you to tune programs by controlling the placement of memory in NUMA systems.	http://linux.softpedia.com/get/Utilities/numactl-8223.shtml
<u> </u>		OpenCL™ (Open Computing Language) is a low-level API for heterogeneous computing that runs on CUDA-nowered GPUs. Using the OpenCL API	· · · · · · · · · · · · · · · · · · ·
327	opencl	developers can launch compute kernels written using a limited subset of the C programming language on a GPU.	https://developer.nvidia.com/opencl
		The Open System for Earthquake Engineering Simulation (OpenSees) is a software framework for simulating the seismic response of structural and	
328	opensees	geotechnical systems. OpenSees has been developed as the computational platform for research in performance-based earthquake engineering	http://opensees.berkeley.edu/OpenSees/home/about.ohp
320	optseg2	ontsen2 is a tool for automatically scheduling events for ranid-resentation event-related (RPER) MRI eventments	http://sufer.nmr.mdb.barvard.edu/optiseo/
330	ovito	OVITO is a scientific visualization and analysis software for atomistic simulation data	http://www.ovito.org/
550		The PCRE library is a set of functions that implement regular expression pattern matching using the same syntax and semantice as Parl 5. PCRE has its	
		own native API as well as a set of wrapper functions that correspond to the POSIX regular expression API. The PCRF library is free even for building	
331	ncre		http://www.pcre.org/
331	000	PHREFECC version 3 is a computer program written in the C and C++ programming languages that is designed to parform a wide version of aqueous	ndavi u u borovo di
332	phreeac	nenchanical calculations	http://www.htr.cr.usas.gov/projects/GWC.coupled/phreegc/
552		PI INK is a free open-source whole genome association analysis toolset, designed to perform a range of basic, large-scale analysis in a computationally	unter un constructed automotion principlicades
332	plink	afficient manner	http://pngu.mgb.barvard.edu/~purcell/plipk/
224	pourou	consolina mannor.	http://www.pourou.vo/
335	promay	a day adong program when generates images non a text based source description, and is available for a variety of computer pidlionits.	Ing.//www.portay.org/
333	promax	PMani is a CIII for PWoof based prearange from Quentum ESPRESSO integrated quite of order for electronic structure selectedates and metacida	
226	pwgui	r wyu is a contor r wso based programs nom quantum-ESPRESSO megrated suite of codes for electronic structure calculations and materials	http://www.k2.ijc.ci/kokoli/owoui/
336	pwgui	moderning at the manoscale	http://www-ko.jp.s/kokal/pwgu/
007		pyr can is a ryuron non-end to the Open-OAM (Open Source CFD Tooloox). It introduces interactivity into Open-OAM, simplifies connection with third-	http://www.facenewillia.com/facenewillia.com/facenewillia.com/facenewillia.com/facenewillia.com/facenewillia.com
33/	pyroan	party functionality and subamilines design of custom user solvers	Intp://openicaniwiki.net/index.pnp/Control/Pyr-cam
338	рупо	r ymou is a user-sponsored morecular visualization system on an open-source foundation.	nup.//www.pymor.org/
		unui computes the convex hull, Delaunay trangulation, voronoi diagram, halfspace intersection about a point, furthest-site Delaunay triangulation, and	
		Turtnest-site Voronoi diagram. It runs in 2-d, 3-d, 4-d, and higher dimensions. It implements the Quickhull algorithm for computing the convex hull. Qhull	
339	qhull	nandies roundoff errors from floating point arithmetic. It can approximate a convex hull.	http://www.qhull.org/
340	qiime	is an open source software package for comparison and analysis of microbial communities, primarily based on high-throughput amplicon sequencing data	http://qiime.org/
341	realvnc	VNC enables you to remotely access and control your devices wherever you are in the world, whenever you need to	https://www.realvnc.com/
342	rnamotif	namotif program [1] searches a database for RNA sequences that match a "motif" describing secondary structure interactions.	http://casegroup.rutgers.edu/casegr-sh-2.5.html

	A	S S	T
		is a problem solving environment or "computational workbench" in which a user selects software modules that can be connected in a visual programing	
242	opinup	is a problem solving characteria in the layer workdown on the manufacture and solver solvers solvers and can be connected in a visual programming	http://www.coi.utok.colu/cike.coftuoro/coirup.html
343	sciluit	environment to create a nigh level workidw for experimentation	http://www.sci.utan.edu/cibc-software/scirun.html
344	sedumi	sortware for optimization over symmetric cones	http://sedumi.ie.lenigh.edu/
345	seg3D	Seg3D is a free volume segmentation and processing tool developed by the NIH Center	http://www.sci.utah.edu/cibc-software/seg3d.html
		SIESTA is both a method and its computer program implementation, to perform efficient electronic structure calculations and ab initio molecular dynamics	
		simulations of molecules and solids. SIESTA's efficiency stems from the use of strictly localized basis sets and from the implementation of linear-scaling	
		algorithms which can be applied to suitable systems. A very important feature of the code is that its accuracy and cost can be tuned in a wide range from	
246	ciosta	uick exploratory calculations to highly accurate simulations matching the quality of other approaches, such as plane-waye and all-electron methods	http://dopartmonts.icmah.os/loom/sigsta/
340	SIESIA	quick exploratory executions to regime accurate simulations matching the quarky of other approaches, other as practices and an execution metalogic exploratory execution and an execution of the quarky of other approaches, doer as practices and an execution metalogic execution and an execution of the quarky of other approaches, doer as practices and an execution of the quarky of other approaches, doer as practices and an execution of the quarky of other approaches, doer as practices and an execution of the quarky of other approaches, doer as practices and an execution of the quarky of other approaches, doer as practices and an execution of the quarky of other approaches, doer as practices and an execution of the quarky of other approaches, doer as practices and an execution of the quarky of other approaches, doer as practices and an execution of the quarky of other approaches, doer as practices and an execution of the quarky of other approaches, doer as practices and an execution of the quarky of other approaches, doer as practices and an execution of the quarky of other approaches, doer as practices and an execution of the quarky of other approaches, doer as practices and an execution of the quarky of other approaches, doer as practices and an execution of the quarky of other approaches, doer as practices and an execution of the quarky of other approaches, doer as practices and an execution of the quarky of other approaches, doer as practices and an execution of the quarky of other approaches, doer as practices and an execution of the quarky of other approaches, doer as practices and an execution of the quarky of other approaches, doer as practices and an execution of the quarky of other approaches, doer as practices and an execution of the quarky of other approaches, doer as practices and an execution of the quarky of other approaches, doer as practices and an execution of the quarky of other approaches, doer as practices and an execution of the quarky of	nitp://departments.icmab.es/ieem/siesta/
0.47		program package for the modeling of atomic resolution structures into low-resolution density maps e.g. non-electron microscopy, tomography, or small	
347	Situs	angle X-ray scattering	http://situs.biomachina.org/
348	socnetv	Social Networks Visualizer (SocNetV) is a cross-platform, user-triendly tool for the analysis and visualization of Social Networks	http://socnetv.sourceforge.net/
349	spider		
350	spm.python		
351	STRanger		
352	teem		
353	Terachem	TeraChem is general numose quantum chemistry software designed to run on NVIDIA GPU architectures under a 64-bit Linux operating system	http://www.petachem.com/index.html
254	v cim	V Sim visualizas atomis structures such as autoritade analysis to rain on the many file formate, binard at site attraction of the second s	http://www.petaenemiconfinitionality
055	v_siin		nitp://www-drintc.cea.n/t_3m/v_3m/
300	veda		
		VirtualGL is an open source toolkit that gives any Unix or Linux remote display software the ability to run OpenGL applications with full 3D hardware	
356	virtualgl	acceleration	http://www.virtualgl.org/
357	wien2k	perform electronic structure calculations of solids using density functional theory (DFT)	http://www.wien2k.at/
358	wordnet	WordNet® is a large lexical database of English	http://wordnet.princeton.edu/wordnet/download/
359	wxDragon	?	http://www.wxdragon.de/
1		XEditd Release 7 (XE7) is the market's most modern 3D electromagnetic simulation software for EDTD-based modeling and simulation. It's tremendously	
260	XEdtd	fact by laveraning NUDIA's most advanced CLIDA anabled GPUIs And it's apply to use an analized an advanced to realized and user advanced of the second	http://www.remcom.com/vf7
300	Ar did	Place by reverging revisions most advanced or Don-reliable or Donating is easy to use and engineered to repricate real-world processes.	ntq.//www.acmcont.com/xr/
	t t	In A moveming anguage to a variate inducing and solution of convex and nonconvex optimization problems, it is implemented as a free (as in a choice) task as (ASTA)	
361	yaimip	no charge) toolbox for MATLAB.	http://users.isy.iiu.se/johani/yalmip/
362			
363	hadoop	Hadoop is an open source data processing framework.	http://hadoop.apache.org/
364	hbase	Hbase is a part of Hadoop, an open source data processing framework.	http://hbase.apache.org/
365	hive	Hive is a part of Hadoop, an open source data processing framework.	http://hive.apache.org/
366	moah	loade the Moah scheduling command environment	http://www.adaptivecomputing.com/
267	nia	Da is a near scheduling command environment	http://www.adabivecompany.com/
307	pig	rig is an open source data processing tool.	
308	sqoop	Squup is a part of haddop; an open source data processing namework.	nttp://sqoop.apacne.org/
369	apache subversion		
370	acmi	ACML is the AMD Core Math Library, which contains functions from BLAS and LAPACK, among other things, optimized for the AMD CPUs	http://developer.amd.com/tools-and-sdks/cpu-development/amd-core-math-library-acml/
		The Enthought Python Distribution provides scientists with a comprehensive set of tools to perform rigorous data analysis and visualization. Python,	
		distinguished by its flexibility, coherence, and ease-of-use, is rapidly becoming the programming language of choice for researchers worldwide. EPD	
		extends this capacity with a powerful collection of Python libraries to enable interactive technical computing and cross-platform rapid application	
371	end	development. This module also includes overc	http://www.enthought.com/
272	floxim	The Elevine is a common floating licease manager commonly used by applications	http://www.flovpooffuero.com/producte/flovpot.publichor.htm
312	liexiii	The reading is a common moding inclusion manager commonly used by applications	nitp://www.nexerasonware.com/products/nexitet/publisher.ntm
0.70		Users should use grottan rather than Gost Gost is a stable, production Portrain So complete available for multiple opti architectures and operating systems.	
373	g95	Innovations and optimizations continue to be worked on. Parts of the F2003 and F2008 standards have been implemented in g95	nttp://g95.org/
		HDF5 for Python (h5py) is a general-purpose Python interface to the Hierarchical Data Format library, version 5. HDF5 is a versatile, mature scientific	
374	h5py	software library designed for the fast, flexible storage of enormous amounts of data.	http://code.google.com/p/h5py/
375	imsl	?	
		MPI for Python (mpi4py) provides bindings of the Message Passing Interface (MPI) standard for the Python programming language. allowing any Python	
376	mpi4py	program to exploit multiple processors.	http://mpi4py.scipy.org/
377	nag	NAGWARE – 195/190 fortran compiler	
270	naglib	NAC has over 1.450 tried and toted routines that are both flavible and portable	http://www.nag.co.dk/
510	l	Derfe vice or automatically analyzes the performance of province and suggests estimizations to allowinte the identified battlengate. It is intended to make	mp.// mm.mag.co.ad/humenoriumenori_initanea.aap
1	1	prometer automatically analyzes the performance or programs and suggests optimizations to alleviate the identified bottlenecks. It is intended to make	
4		and a many second second the second state of the second se	
	perrexpert	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it	
379	let i	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled down data sets and resource configurations to obtain accurate analyses.	http://www.tacc.utexas.edu/perfexpert/
379 380	Rhadoop	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework.	http://www.tacc.utexas.edu/perfexpert/ https://github.com/RevolutionAnalytics/RHadoop/wiki
379 380	Rhadoop	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework. CFOUR (Coupled-Cluster techniques for Computational Chemistry) is a program package for performing high-level quantum mechanical calculations on	http://www.tacc.utexas.edu/perfexpert/ https://github.com/RevolutionAnalytics/RHadoop/wiki
379 380 <u>38</u> 1	Rhadoop cfour	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework. CFOUR (Coupled-Cluster techniques for Computational Chemistry) is a program package for performing high-level quantum mechanical calculations on atoms and molecules.	http://www.tacc.utexas.edu/perfexpert/ https://github.com/RevolutionAnalytics/RHadoop/wiki http://www.cfour.de/
379 380 381	Rhadoop cfour	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework. CFOUR (Coupled-Cluster techniques for Computational Chemistry) is a program package for performing high-level quantum mechanical calculations on atoms and molecules. Chombo provides a set of tools for implementing finite difference methods for the solution of partial differential equations on block-structured adaptively.	http://www.tacc.utexas.edu/perfexpert/ https://github.com/RevolutionAnalytics/RHadoop/wiki http://www.cfour.de/
379 380 381	Rhadoop cfour	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework. CFOUR (Coupled-Cluster techniques for Computational Chemistry) is a program package for performing high-level quantum mechanical calculations on atoms and molecules. Chombo provides a set of tools for implementing finite difference methods for the solution of partial differential equations on block-structured adaptively refined rectangular grids. Both elliptic and time-dependent modules are included. Chombo supports calculations in commerce with hoth	http://www.tacc.utexas.edu/perfexpert/ https://github.com/RevolutionAnalytics/RHadoop/wiki http://www.cfour.de/
379 380 381	Rhadoop cfour	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework. CFOUR (Coupled-Cluster techniques for Computational Chemistry) is a program package for performing high-level quantum mechanical calculations on atoms and molecules. Chombo provides a set of tools for implementing finite difference methods for the solution of partial differential equations on block-structured adaptively refined rectangular grids. Both elliptic and time-dependent modules are included. Chombo supports calculations in complex geometries with both embedded boundaries and mapped drids, and Chombo supports calculations. Most narialle inlatforms are supnorted and cress-inlatform self.	http://www.tacc.utexas.edu/oerfexpert/ https://github.com/RevolutionAnalytics/RHadoop/wiki http://www.cfour.de/
379 380 381	Rhadoop cfour	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework. CFOUR (Coupled-Cluster techniques for Computational Chemistry) is a program package for performing high-level quantum mechanical calculations on atoms and molecules. Chombo provides a set of tools for implementing finite difference methods for the solution of partial differential equations on block-structured adaptively refined rectangular grids. Both elliptic and time-dependent modules are included. Chombo supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self- describing if formats are included.	http://www.tacc.utexas.edu/perfexpert/ https://github.com/RevolutionAnalytics/RHadoop/wiki http://www.cfour.de/
379 380 381	Rhadoop cfour chombo	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework. CFOUR (Coupled-Cluster techniques for Computational Chemistry) is a program package for performing high-level quantum mechanical calculations on atoms and molecules. Chombo provides a set of tools for implementing finite difference methods for the solution of partial differential equations on block-structured adaptively refined rectangular grids. Both elliptic and time-dependent modules are included. Chombo supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self- describing file formats are included.	http://www.tacc.utexas.edu/perfexpert/ https://github.com/RevolutionAnalytics/RHadoop/wiki http://www.cfour.de/ https://commons.lbl.gov/display/chombo/Chombo+++Software+for+Adaptive+Solutions+of+Partial+Differential+Equations
379 380 381 382 382 383	Rhadoop cfour chombo dakota	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework. CFOUR (Coupled-Cluster techniques for Computational Chemistry) is a program package for performing high-level quantum mechanical calculations on atoms and molecules. Chombo provides a set of tools for implementing finite difference methods for the solution of partial differential equations on block-structured adaptively refined rectangular grids. Both elliptic and time-dependent modules are included. Chombo supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self- describing file formats are included. Dakota — A Mutilitee Parallel Object-Oriented Framework for design optimization	http://www.tacc.utexas.edu/perfexpert/ http://github.com/RevolutionAnalytics/RHadoop/wiki http://www.cfour.de/ https://commons.ibl.gov/display/chombo/Chombo+-+Software+for+Adaptive+Solutions+of+Partial+Differential+Equations http://dakota.sandia.gov/
379 380 381 382 383 383	Rhadoop cfour chombo dakota globus	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework. CFOUR (Coupled-Cluster techniques for Computational Chemistry) is a program package for performing high-level quantum mechanical calculations on atoms and molecules. Chombo provides a set of tools for implementing finite difference methods for the solution of partial differential equations on block-structured adaptively refined rectangular grids. Both elliptic and time-dependent modules are included. Chombo supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self- describing file formats are included. Daktai — A Multilevel Parallel Object-Oriented Framework for design optimization Globus provides the GSI/MyProxy in addition to GridFTP cli tools	http://www.tacc.utexas.edu/perfexpert/ https://github.com/RevolutionAnalytics/RHadoop/wiki http://www.cfour.de/ https://commons.lbl.gov/display/chombo/Chombo+-+Software+for+Adaptive+Solutions+of+Partial+Differential+Equations http://dakota.sandia.gov/ http://dakota.sandia.gov/
379 380 381 382 383 384 385	Rhadoop cfour chombo dakota globus gti	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework. CFOUR (Coupled-Cluster techniques for Computational Chemistry) is a program package for performing high-level quantum mechanical calculations on atoms and molecules. Chombo provides a set of tools for implementing finite difference methods for the solution of partial differential equations on block-structured adaptively refined rectangular grids. Both elliptic and time-dependent modules are included. Chombo supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self- describing file formats are included. Chombo also Supports particle methods. Most parallel platforms are supported, and cross-platform self- describing file formats are included. Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self- describing file formats are included. Diset-Oriented Framework for design optimization Globus provides the GSIMyProxy in addition to clo, capable of integrated simulations of the total engine and powertrain system	http://www.tacc.utexas.edu/perfexpert/ https://github.com/RevolutionAnalytics/RHadoop/wiki http://www.cfour.de/ https://commons.ibl.gov/display/chombo/Chombo+-+Software+for+Adaptive+Solutions+of+Partial+Differential+Equations http://dakota.sandia.gov/ http://www.globus.org/ http://www.globus.org/
379 380 381 382 383 384 385	Rhadoop cfour chombo dakota globus gti	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework. CFOUR (Coupled-Cluster techniques for Computational Chemistry) is a program package for performing high-level quantum mechanical calculations on atoms and molecules. Chombo provides a set of tools for implementing finite difference methods for the solution of partial differential equations on block-structured adaptively refined rectangular grids. Both elliptic and time-dependent modules are included. Chombo supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self- describing file formats are included. Daktos — A Multileve Plarallel Object-Oriented Framework for design optimization Globus provides the GSI/MyProxy in addition to GridFTP cli tools GT-SUITE offers the only true virtual engine/powertrain tool, capable of integrated simulations of the total engine and powertrain system	http://www.tacc.utexas.edu/perfexpert/ https://github.com/RevolutionAnalytics/RHadoop/wiki http://www.cfour.de/ https://commons.tbl.gov/display/chombo/Chombo+-+Software+for+Adaptive+Solutions+of+Partial+Differential+Equations http://dakota.sandia.gov/ http://www.globus.org/ http://www.globus.org/
379 380 381 382 383 384 385	Rhadoop cfour chombo dakota globus gti	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework. CFOUR (Coupled-Cluster techniques for Computational Chemistry) is a program package for performing high-level quantum mechanical calculations on atoms and molecules. Chombo provides a set of tools for implementing finite difference methods for the solution of partial differential equations on block-structured adaptively refined rectangular grids. Bott elliptic and time-dependent modules are included. Chombo supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self- describing file formats are included. Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self- describing file formats are included. Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self- describing file formats are included. Dakota — A Mutilievel Parallel Object-Oriented Framework for design optimization Globus provides the SliMyProxy in addition to GridFTP cli tools GT-SUITE offers the only true virtual engine/powertrain tool, capable of integrated simulations of the total engine and powertrain system HPCToolkit is an integrated suite of tools for measurement and analysis of program performance on computers ranging from multicore desktop systems to	http://www.tacc.utexas.edu/perfexpert/ https://github.com/RevolutionAnalytics/RHadoop/wiki http://www.cfour.de/ https://commons.lbl.gov/display/chombo/Chombo+-+Software+for+Adaptive+Solutions+of+Partial+Differential+Equations http://dakota.sandia.gov/ http://www.gtisoft.com/
379 380 381 382 383 384 385	Rhadoop cfour chombo dakota globus gti	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework. CFOUR (Coupled-Cluster techniques for Computational Chemistry) is a program package for performing high-level quantum mechanical calculations on atoms and molecules. Chombo provides a set of tools for implementing finite difference methods for the solution of partial differential equations on block-structured adaptively refined rectangular grids. Both elliptic and time-dependent modules are included. Chombo supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self- describing file formats are included. Dakta — A Multilevel Parallel Object-Oriented Framework for design optimization Globus provides the GSI/MyProxy in addition to GridFTP cli tools GT-SUITE offers the only true virtual engine/powertrain tool, capable of integrated simulations of the total engine and powertrain system HPCToolkit is an integrated suite of tools for measurement and analysis of program performance on computers ranging from multicore desktop systems to the nation's largest supercomputers. By using statistical sampling of timers and hardware performance counters, HPCToolkit collects accurate	http://www.tacc.utexas.edu/perfexpert/ https://github.com/RevolutionAnalytics/RHadoop/wiki http://www.cfour.de/ https://commons.ibl.gov/display/chombo/Chombo++Software+for+Adaptive+Solutions+of+Partial+Differential+Equations http://dakota.sandia.gov/ http://www.globus.org/ http://www.globus.org/
379 380 381 382 383 384 385 386	Rhadoop cfour chombo dakota globus gti hpctoolkit	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework. CFOUR (Coupled-Cluster techniques for Computational Chemistry) is a program package for performing high-level quantum mechanical calculations on atoms and molecules. Chombo provides a set of tools for implementing finite difference methods for the solution of partial differential equations on block-structured adaptively refined rectangular grids. Both elliptic and time-dependent modules are included. Chombo supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self- describing file formats are included. Dakota — A Mutilievel Parallel Object-Oriented Framework for design optimization Globus provides the GSI/MyProxy in addition to GridFTP cli tools GT-SUITE offers the only true virtual engine/powertrain tool, capable of integrated simulations of the total engine and powertrain system HPCToolkit is an integrated suite of tools for measurement and analysis of program performance on computers ranging from multicore desktop systems to the nation's largest supercomputers. By using statistical sampling of timers and hardware performance counters, HPCToolkit collects accurate measurements of a program's work, resource consumption, and intefficiency and attributes them to the full calino context in which the view counters in which the view counter super context of a program's work, resource consumption, and integricery and attributes them to the full calino context in which the view counter super context of a program's work, resource consumption, and integricery and attributes them to the full calino context on work in work.	http://www.tacc.utexas.edu/perfexpert/ https://github.com/RevolutionAnalytics/RHadoop/wiki http://www.cfour.de/ http://commons.lbl.gov/display/chombo/Chombo+-+Software+for+Adaptive+Solutions+of+Partial+Differential+Equations http://dakota.sandia.gov/ http://www.gtisoft.com/ http://
379 380 381 382 383 384 385 386 387	Rhadoop cfour chombo dakota globus gli hpctoolkit monaodh	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework. CFOUR (Coupled-Cluster techniques for Computational Chemistry) is a program package for performing high-level quantum mechanical calculations on atoms and molecules. Chombo provides a set of tools for implementing finite difference methods for the solution of partial differential equations on block-structured adaptively refined rectangular grids. Both elliptic and time-dependent modules are included. Chombo supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self- describing file formats are included. Dakta – A Nutlikevel Parallel Object-Oriented Framework for design optimization Globus provides the GSI/MyProxy in addition to GridFTP cli tools GT-SUITE offers the only true virtual engine/powertrain tool, capable of integrated simulations of the total engine and powertrain system HPCToolkit is an integrated suite of tools for measurement and analysis of program performance on computers ranging from multicore desktop systems to the nation's largest supercomputers. By using statistical sampling of timers and hardware performance counters, HPCToolkit collects accurate measurements of a program's work, resource consumption, and inefficiency and attributes them to the full calling context in which they occur. 7	http://www.tacc.utexas.edu/perfexpert/ https://github.com/RevolutionAnalytics/RHadoop/wiki http://www.cfour.de/ https://commons.ibl.gov/display/chombo/Chombo+-+Software+for+Adaptive+Solutions+of+Partial+Differential+Equations http://dakota.sandia.gov/ http://www.globus.org/ http://www.globus.org/
379 380 381 382 383 384 385 386 387	Rhadoop cfour chombo dakota globus gti hpctoolkit mongodb	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework. CFOUR (Coupled-Cluster techniques for Computational Chemistry) is a program package for performing high-level quantum mechanical calculations on atoms and molecules. Chombo provides a set of tools for implementing finite difference methods for the solution of partial differential equations on block-structured adaptively refined rectangular grids. Both elliptic and time-dependent modules are included. Chombo supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self- describing file formats are included. Dakota — A Mutilievel Parallel Object-Oriented Framework for design optimization Globus provides the GSI/MyProxy in addition to GridFTP di tools GT-SUITE offers the only true virtual engine/powertrain tool, capable of integrated simulations of the total engine and powertrain system HPCToolkit is an integrated suite of tools for measurement and analysis of program performance on computers ranging from multicore desktop systems to the nation's largest supercomputers. By using statistical sampling of timers and hardware performance counters, HPCToolkit collects accurate measurements of a program's work, resource consumption, and inefficiency and attributes them to the full calling context in which they occur. 7	http://www.tacc.utexas.edu/perfexpert/ https://github.com/RevolutionAnalytics/RHadoop/wiki http://www.cfour.de/ https://commons.lbl.gov/display/chombo/Chombo++Software+for+Adaptive+Solutions+of+Partial+Differential+Equations http://akota.sandia.gov/ http://www.glisoft.com/ http://htpctoolkit.org/
379 380 381 382 383 384 385 386 387	Rhadoop cfour chombo dakota globus gti hpctoolkit mongodb	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework. CFOUR (Coupled-Cluster techniques for Computational Chemistry) is a program package for performing high-level quantum mechanical calculations on atoms and molecules. Chombo provides a set of tools for implementing finite difference methods for the solution of partial differential equations on block-structured adaptively refined rectangular grids. Both elliptic and time-dependent modules are included. Chombo supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self- describing file formatis are included. Dakta — A Nutlikevel Parallel Object-Oriented Framework for design optimization Globus provides the GSI/MyProxy in addition to GridFTP cli tools GT-SUITE offers the only true virtual engine/powertrain tool, capable of integrated simulations of the total engine and powertrain system HPCToolkit is an integrated suite of tools for measurement and analysis of program performance on computers ranging from multicore desktop systems to the nation's largest supercomputers. By using statistical sampling of timers and hardware performance counters, HPCToolkit collects accurate measurements of a program's work, resource consumption, and inefficiency and attributes them to the full calling context in which they occur. ?	http://www.tacc.utexas.edu/perfexpert/ https://github.com/RevolutionAnalytics/RHadoop/wiki http://www.cfour.de/ http://commons.ibl.gov/display/chombo/Chombo+-+Software+for+Adaptive+Solutions+of+Partial+Differential+Equations http://dakota.sandia.gov/ http://www.globus.org/ http://www.globus.org/
379 380 381 382 383 384 385 385 386 387	Rhadoop cfour chombo dakota globus gti hpctoolkit mongodb	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework. CFOUR (Coupled-Cluster techniques for Computational Chemistry) is a program package for performing high-level quantum mechanical calculations on atoms and molecules. Chombo provides a set of tools for implementing finite difference methods for the solution of partial differential equations on block-structured adaptively refined rectangular grids. Both elliptic and time-dependent modules are included. Chombo supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self-describing file formats are included. Chombo also supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self-describing file formats are included. Chombo data — A Mutilievel Parallel Object-Oriented Framework for design optimization Giobus provides the GSI/MyProxy in addition to GridFTP cli tools GT-SUITE offers the only true virtual engine/powertrain tool, capable of integrated simulations of the total engine and powertrain system HPCToolkit is an integrated suite of tools for measurement and analysis of program performance on computers ranging from multicore desktop systems to the nation's largest supercomputers. By using statistical sampling of timers and hardware performance counters, HPCToolkit collects accurate measurements of a program swork, resource consumption, and inefficiency and attributes them to the full calling context in which they occur. 7 This program computes definite-frequency eigenstates (harmonic modes) of Maxwell's e	http://www.tacc.utexas.edu/perfexpert/ https://github.com/RevolutionAnalytics/RHadoop/wiki http://www.cfour.de/ https://commons.lbl.gov/display/chombo/Chombo+++Software+for+Adaptive+Solutions+of+Partial+Differential+Equations http://dakota.sandia.gov/ http://www.glisoft.com/ http://hpctoolkit.org/
379 380 381 382 383 383 385 385 386 387	Rhadoop ctour chombo dakota globus gli hpctoolkit mongodb	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework. CFOUR (Coupled-Cluster techniques for Computational Chemistry) is a program package for performing high-level quantum mechanical calculations on atoms and molecules. Chombo provides a set of tools for implementing finite difference methods for the solution of partial differential equations on block-structured adaptively refined rectangular grids. Both elliptic and time-dependent modules are included. Chombo supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self- describing file formatis are included. Daktat — A Multilevel Parallel Object-Oriented Framework for design optimization GT-SUITE offers the only true virtual engine/powertrain tool, capable of integrated simulations of the total engine and powertrain system HPCToolkit is an integrated suite of tools for measurement and analysis of program performance on computers ranging from multicore desktop systems to the nation's largest supercomputers. By using statistical sampling of timers and hardware performance counters, HPCToolkit collects accurate measurements of a program's work, resource consumption, and inefficiency and attributes them to the full calling context in which they occur. 7 This program computes definite-frequency eigenstates (harmonic modes) of Maxwell's equations in periodic dielectric structures for arbitrary wavevectors, also applicable to mony other problems in optics, such as awaeguides and resonator systems. (For example, it can solve for the modes of waveguides with slow applicable to mony other problems in optics, such as waveguides and resonator systems	http://www.tacc.utexas.edu/perfexpert/ https://github.com/RevolutionAnalytics/RHadoop/wiki http://www.cfour.de/ http://commons.ibl.gov/display/chombo/Chombo+-+Software+for+Adaptive+Solutions+of+Partial+Differential+Equations http://dakota.sandia.gov/ http://www.globus.org/ http://htpctoolkit.org/
379 380 381 382 383 384 385 386 387 388	Rhadoop cfour chombo dakota globus gti hpctoolkit mongodb mpb	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework. CFOUR (Coupled-Cluster techniques for Computational Chemistry) is a program package for performing high-level quantum mechanical calculations on atoms and molecules. Chombo provides a set of tools for implementing finite difference methods for the solution of partial differential equations on block-structured adaptively refined rectangular grids. Both elliptic and time-dependent modules are included. Chombo supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self-describing file formats are included. Chombo also supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self-describing file formats are included. Chombo also supports parallel platforms are supported, and cross-platform self-describing file formats are included. The provide is the GSI/MyProxy in addition to GridFTP cli tools GT-SUITE offers the only true virtual engine/powertrain tool, capable of integrated simulations of the total engine and powertrain system HPCToolkit is an integrated suite of tools for measurement and analysis of program performance on computers ranging from multicore desktop systems to the nation's largest supercomputers. By using statistical sampling of timers and hardware performance counters, HPCToolkit collects accurate measurements of a program's work, resource consumption, and inefficiency and attributes them to the full calling context in which they occur. 7. This program computes definite-frequency eigenstates (har	http://www.tacc.utexas.edu/perfexpert/ https://github.com/RevolutionAnalytics/RHadoop/wiki http://www.cfour.de/ https://commons.lbl.gov/display/chombo/Chombo+-+Software+for+Adaptive+Solutions+of+Partial+Differential+Equations http://akota.sandia.gov/ http://www.gitsoft.com/ http://hpctoolkit.org/
379 380 381 382 383 384 385 386 387 388 388 388	Rhadoop ctour chombo dakota globus gli hpctoolkit mongodb mpb mpcci	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework. CFOUR (Coupled-Cluster techniques for Computational Chemistry) is a program package for performing high-level quantum mechanical calculations on atoms and molecules. Chombo provides a set of tools for implementing finite difference methods for the solution of partial differential equations on block-structured adaptively refined rectangular grids. Both elliptic and time-dependent modules are included. Chombo supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self- describing file formatis are included. Daktat — A Multiblevel Parallel Object-Oriented Framework for design optimization Globus provides the GSI/MyProxy in addition to GridFTP cli tools GT-SUITE offers the only true virtual engine/powertrain tool, capable of integrated simulations of the total engine and powertrain system HPCToolkit is an integrated suite of tools for measurement and analysis of program performance on computers ranging from multicore desktop systems to the nation's largest supercomputers. By using statistical sampling of timers and hardware performance counters, HPCToolkit collects accurate measurements of a program's work, resource consumption, and inefficiency and attributes them to the full calling context in which they occur. 7 This program computes definite-frequency eigenstates (harmonic modes) of Maxwell's equations in periodic dielectric structures for arbitrary wavevectors. using fully-vectorial and three-demensional methods. It is especially designed for the study of photonic carysta (a.k.a. photonic band-gap materilab), but is also applicable t	http://www.tacc.utexas.edu/perfexpert/ https://github.com/RevolutionAnalytics/RHadoop/wiki http://www.cfour.de/ http://commons.ibl.gov/display/chombo/Chombo+-+Software+for+Adaptive+Solutions+of+Partial+Differential+Equations http://dakota.sandia.gov/ http://dakota.sandia.gov/ http://www.globus.org/ http://htpctoolkit.org/ http://htpctoolkit.org/
379 380 381 382 383 384 385 386 387 388 389	Rhadoop cfour chombo dakota globus gli hpctoolkit mongodb mpb mpcci	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework. CFOUR (Coupled-Cluster techniques for Computational Chemistry) is a program package for performing high-level quantum mechanical calculations on atoms and molecules. Chombo provides a set of tools for implementing finite difference methods for the solution of partial differential equations on block-structured adaptively refined rectangular grids. Both elliptic and time-dependent modules are included. Chombo supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self-describing file formats are included. Chombo also supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self-describing file formats are included. Chombo also supports parallel platforms are supported, and cross-platform self-describing file formats are included. The complex second set of SIMP Proxy in addition to GridFTP cli tools GT-SUITE offers the only true virtual engine/powertrain tool, capable of integrated simulations of the total engine and powertrain system HPCToolkit is an integrated suite of tools for measurement and analysis of program performance on computers ranging from multicore desktop systems to the nation's largest supercomputers. By using statistical sampling of timers and hardware performance counters, HPCToolkit collects accurate measurements of a program 'sowrk, resource consumption, and inefficiency and attributes them to the full calling context in which they occur. 7 This program computes definite-frequency eigenstates	http://www.tacc.utexas.edu/perfexpert/ https://github.com/RevolutionAnalytics/RHadoop/wiki http://www.cfour.de/ https://commons.lbl.gov/display/chombo/Chombo+-+Software+for+Adaptive+Solutions+of+Partial+Differential+Equations http://akota.sandia.gov/ http://www.gtisoft.com/ http://hpctoolkit.org/ http://hpctoolkit.org/
379 380 381 383 383 384 385 386 387 388 389 390	Rhadoop cfour chombo dakota globus gli hpctoolkit mongodb mpb mpcci ncl	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework. CFOUR (Coupled-Cluster techniques for Computational Chemistry) is a program package for performing high-level quantum mechanical calculations on atoms and molecules. Chombo provides a set of tools for implementing finite difference methods for the solution of partial differential equations on block-structured adaptively refined rectangular grids. Both elliptic and time-dependent modules are included. Chombo supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self- describing file formats are included. Dakota — A Multilevel Parallel Object-Oriented Framework for design optimization Globus provides the GSI/MyProxy in addition to GrifFTP cli tools GT-SUITE offers the only true virtual engine/powertrain tool, capable of integrated simulations of the total engine and powertrain system HPCToolkit is an integrated suite of tools for measurement and analysis of program performance on computers ranging from multicore desktop systems to the nation's largest supercomputers. By using statistical sampling of timers and hardware performance counters, HPCToolkit collects accurate measurements of a program's work, resource consumption, and inefficiency and attributes them to the full calling context in which they occur. 7 This program computes definite-frequency eigenstates (harmonic modes) of Maxwell's equations in periodic dielectric structures for arbitrary wavevectors, and pluty-accoratic and three-dimensional methods. It is especially designed for the study of photonic cortsal (ask a. phonic band-gap materilas), but is also apolicable to many	http://www.acc.utexas.edu/perfexpert/ https://github.com/RevolutionAnalytics/RHadoop/wiki http://www.cfour.de/ http://commons.lbl.gov/display/chombo/Chombo+-+Software+for+Adaptive+Solutions+of+Partial+Differential+Equations http://dakota.sandia.gov/ http://dakota.sandia.gov/ http://www.globus.org/ http://hpctoolkit.org/ http://hpctoolkit.org/ http://hpctoolkit.org/ http://ab-initio.mit.edu/wiki/index.php/MIT_Photonic_Bands http://www.mpcci.de/multiphysics-engineering.html http://www.ncl.ucar.edu/
379 380 381 382 383 384 385 386 387 388 389 389 390	Rhadoop cfour chombo dakota globus gli hpctoolkit mongodb mpc mpc ncl	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework. CFOUR (Coupled-Cluster techniques for Computational Chemistry) is a program package for performing high-level quantum mechanical calculations on atoms and molecules. Chombo provides a set of tools for implementing finite difference methods for the solution of partial differential equations on block-structured adaptively refined rectangular grids. Both elliptic and time-dependent modules are included. Chombo supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self-describing file formats are included. Chombo also supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self-describing file formats are included. Chombo also supports parallel platforms are supported, and cross-platform self-describing file formats are included. The complex second setting of the solution of the total engine and powertrain system HPCToolkit is an integrated suite of tools for measurement and analysis of program performance on computers ranging from multicore desktop systems to the nation's largest supercomputers. By using statistical sampling of timers and hardware performance counters, HPCToolkit collects accurate measurements of a program 's work, resource consumption, and inefficiency and attributes them to the full calling context in which they occur. 7 This program computes definite-frequency eigenstates (harmonic modes) of Maxwell's equations in periodic dielectric structures for arabitrary wavevectors, using fully-vectori	http://www.tacc.utexas.edu/perfexpert/ https://github.com/RevolutionAnalytics/RHadoop/wiki http://commons.lbl.gov/display/chombo/Chombo+++Software+for+Adaptive+Solutions+of+Partial+Differential+Equations http://dakota.sandia.gov/ http://www.gtisoft.com/ http://hpctoolkit.org/ http://hpctoolkit.org/ http://ab-initio.mit.edu/wiki/index.php/MIT_Photonic_Bands http://www.mccl.ucar.edu/
379 380 381 382 383 384 385 388 389 388 389 390	Rhadoop cfour chombo dakota globus globus gti hpctoolkit mongodb mpb mpcci ncl	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework. CFOUR (Coupled-Cluster techniques for Computational Chemistry) is a program package for performing high-level quantum mechanical calculations on atoms and molecules. Chombo provides a set of tools for implementing finite difference methods for the solution of partial differential equations on block-structured adaptively refined rectangular grids. Both elliptic and time-dependent modules are included. Chombo supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self- describing file formats are included. Dakota — A Multilevel Parallel Object-Oriented Framework for design optimization Globus provides the GSI/MyProxy in addition to GrifFTP cli tools GT-SUITE offers the only true virtual engine/powertrain tool, capable of integrated simulations of the total engine and powertrain system HPCToolkit is an integrated suite of tools for measurement and analysis of program performance on computers ranging from multicore desktop systems to the nation's largest supercomputers. By using statistical sampling of timers and hardware performance counters, HPCToolkit collects accurate measurements of a program's work, resource consumption, and inefficiency and attributes them to the full calling context in which they occur. 7 This program computes definite-frequency eigenstates (harmonic modes) of Maxwell's equations. The protonic band-gen materials), but is also applicable to many other problems in optics, such as waveguides and resonator systems. (For example, it can solve for the modes of waveguides with arbitrary cross-sections.). Multiphysics Co	http://www.tacc.utexas.edu/perfexpert/ https://github.com/RevolutionAnalytics/RHadoop/wiki http://www.cfour.de/ http://commons.lbl.gov/display/chombo/Chombo+-+Software+for+Adaptive+Solutions+of+Partial+Differential+Equations http://dakota.sandia.gov/ http://dakota.sandia.gov/ http://bpctoolkit.org/ http://hpctoolkit.org/ http://hpctoolkit.org/
379 380 381 382 383 384 385 386 387 388 389 390	Rhadoop cfour chombo dakota globus gli hpctoolkit mongodb mpb mpcci ncl ochonus	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework. CFOUR (Coupled-Cluster techniques for Computational Chemistry) is a program package for performing high-level quantum mechanical calculations on atoms and molecules. Chombo provides a set of tools for implementing finite difference methods for the solution of partial differential equations on block-structured adaptively refined rectangular grids. Both elliptic and time-dependent modules are included. Chombo supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self-describing file formats are included. Chombo also supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self-describing file formats are included. Chombo also supports parallel platforms are supported, and cross-platform self-describing file formats are included. The program parallel object-Oriented Framework for design optimization. Globus provides the GSI/MyProxy in addition to GridFTP cli tools GT-SUITE offers the only true virtual engine/powertrain tool, capable of integrated simulations of the total engine and powertrain system HPCToolkit is an integrated suite of tools for measurement and analysis of program performance on computers ranging from multicore desktop systems to the nation's largest supercomputers. By using statistical sampling of timers and hardware performance counters, HPCToolkit collects accurate measurements of a program 's work, resource consumption, and inefficiency and attributes them to the full calling context in whic	http://www.tacc.utexas.edu/perfexpert/ https://github.com/RevolutionAnalytics/RHadoop/wiki http://www.cfour.de/ https://commons.lbl.gov/display/chombo/Chombo+++Software+for+Adaptive+Solutions+of+Partial+Differential+Equations http://dakota.sandia.gov/ http://dakota.sandia.gov/ http://www.gtisoft.com/ http://hpctoolkit.org/ http://hpctoolkit.org/ http://ab-initio.mit.edu/wiki/index.php/MIT_Photonic_Bands http://www.mpcci.de/multiphysics-engineering.html http://www.ncci.ucar.edu/
379 380 381 382 383 384 385 386 387 388 389 389 390 391	Rhadoop cfour chombo dakota globus globus gti hpctoolkit mongodb mpb mpcci ncl octopus	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework. CFOUR (Coupled-Cluster techniques for Computational Chemistry) is a program package for performing high-level quantum mechanical calculations on atoms and molecules. Chombo provides a set of tools for implementing finite difference methods for the solution of partial differential equations on block-structured adaptively refinder declangular grids. Both elliptic and time-dependent modules are included. Chombo supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self- describing file formats are included. Dakota — A Multilevel Parallel Object-Oriented Framework for design optimization Globus provides the GSI/MyProxy in addition to GrifFTP cli tools GT-SUITE offers the only true virtual engine/powertrain tool, capable of integrated simulations of the total engine and powertrain system HPCToolkit is an integrated suite of tools for measurement and analysis of program performance on computers ranging from multicore desktop systems to the nation's largest supercomputers. By using statistical sampling of timers and hardware performance counters, HPCToolkit collects accurate measurements of a program's work, resource consumption, and inefficiency and attributes frem to the full calling context in which they occur. 7 This program computes definite-frequency eigenstates (harmonic modes) of Maxwell's equations. The protonic band-gen materials), but is also applicable to many other problems in optics, such as waveguides and resonator systems. (For example, it can solve for the modes of waveguides with arbitrary cross-sections.). Multiphysics C	http://www.tacc.utexas.edu/perfexpert/ https://github.com/RevolutionAnalytics/RHadoop/wiki http://www.cfour.de/ http://commons.lbl.gov/display/chombo/Chombo+-+Software+for+Adaptive+Solutions+of+Partial+Differential+Equations http://dakota.sandia.gov/ http://dakota.sandia.gov/ http://www.globus.org/ http://hpctoolkit.org/ http://hpctoolkit.org/ http://hpctoolkit.org/ http://www.mcci.de/multiphysics-engineering.html http://www.ncl.ucar.edu/ http://www.tddft.org/programs/octopus/wiki/index.php/Main_Page
379 380 381 383 384 385 386 387 388 389 390 390	Rhadoop cfour chombo dakota globus gli hpctoolkit mongodb mpb mpcci ncl octopus	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework. CFOUR (Coupled-Cluster techniques for Computational Chemistry) is a program package for performing high-level quantum mechanical calculations on atoms and molecules. Chombo provides a set of tools for implementing finite difference methods for the solution of partial differential equations on block-structured adaptively refined rectangular grids. Both elliptic and time-dependent modules are included. Chombo supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self-describing file formats are included. Chombo also supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self-describing file formats are included. Chombo also supports calculations of the total engine formation are included. Chombo also supports and the self self self self self self self sel	http://www.tacc.utexas.edu/perfexpert/ https://github.com/RevolutionAnalytics/RHadoop/wiki http://www.cfour.de/ https://commons.lbl.gov/display/chombo/Chombo+++Software+for+Adaptive+Solutions+of+Partial+Differential+Equations http://dakota.sandia.gov/ http://dakota.sandia.gov/ http://www.gtisoft.com/ http://hpctoolkit.org/ http://hpctoolkit.org/ http://hpctoolkit.org/ http://www.mccd.de/multiphysics-engineering.html http://www.utddft.org/porgrams/octopus/wiki/index.php/Main_Page
379 380 381 383 384 385 387 388 389 389 390 391 392	Rhadoop cfour chombo dakota globus gli hpctoolkit mongodb mppc mppci ncl octopus obzip	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework. CFOUR (Coupled-Cluster techniques for Computational Chemistry) is a program package for performing high-level quantum mechanical calculations on atoms and molecules. Chombo provides a set of tools for implementing finite difference methods for the solution of partial differential equations on block-structured adaptively refined rectangular grids. Both elliptic and time-dependent modules are included. Chombo supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self- describing file formats are included. Dakota — A Mutilevel Parallel Object-Oriented Framework for design optimization Globus provides the GSI/MyProxy in addition to GridFTP cli tools GT-SUITE offers the only true virtual engine/powertrain tool, capable of integrated simulations of the total engine and powertrain system HPCToolkit is an integrated suite of tools for measurement and analysis of program performance on computers ranging from multicore desktop systems to the nation's largest supercomputers. By using statistical sampling of timers and hardware performance counters, HPCToolkit collects accurate measurements of a program's work, resource consumption, and inefficiency and attributes filem to the full calling context in which they occur. 7 This program computes definite-frequency eigenstates (harmonic modes) of Maxwell's equations. Toe hotonic band-qape materials), but is also applicable to many other problems in optics, such as waveguides and resonator systems. (For example, it can solve for the modes of waveguides with altoitrary cross-sections.) Multiphysics Co	http://www.tacc.utexas.edu/perfexpert/ https://github.com/RevolutionAnalytics/RHadoop/wiki http://www.cfour.de/ http://commons.lbl.gov/display/chombo/Chombo+-+Software+for+Adaptive+Solutions+of+Partial+Differential+Equations http://dakota.sandia.gov/ http://dakota.sandia.gov/ http://www.globus.org/ http://hpctoolkit.org/ http://hpctoolkit.org/ http://hpctoolkit.org/ http://www.mpcci.de/multiphysics-engineering.html http://www.ncl.ucar.edu/ http://www.ncl.ucar.edu/
379 380 381 382 383 384 385 384 385 386 387 388 389 390 391 392 393	Rhadoop cfour chombo dakota globus gli hpctoolkit mongodb mpb mpcci ncl octopus obzip pcp	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework. CFOUR (Coupled-Cluster techniques for Computational Chemistry) is a program package for performing high-level quantum mechanical calculations on atoms and molecules. Chombo provides a set of tools for implementing finite difference methods for the solution of partial differential equations on block-structured adaptively refined rectangular grids. Both elliptic and time-dependent modules are included. Chombo supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self- describing file formats are included. Dakota — A Mutilievel Parallel Object-Oriented Framework for design optimization Globus provides the GSI/MyProxy in addition to GridFTP cli tools GT-SUITE offers the only true virtual engine/powertrain tool, capable of integrated simulations of the total engine and powertrain system HPCToolkit is an integrated suite of tools for measurement and analysis of program performance on computers ranging from multicore desktop systems to the nation's largest supercomputers. By using statistical sampling of timers and hardware performance counters, HPCToolkit collects accurate measurements of a program's work, resource consumption, and inefficiency and attributes them to the full calling context in which they occur. 7 This program computes definite-frequency eigenstates (harmonic modes) of Maxwell's equations in periodic dielectric structures for arbitrary wavevectors, using fully-vectorial and three-dimensional methods. It is especially designed for the study of photonic crystals (a.k.a. photonic band-gap materials), but is also app	http://www.tacc.utexas.edu/perfexpert/ https://github.com/RevolutionAnalytics/RHadoop/wiki https://commons.lbl.gov/display/chombo/Chombo+++Software+for+Adaptive+Solutions+of+Partial+Differential+Equations http://dakota.sandia.gov/ http://dakota.sandia.gov/ http://www.globus.org/ http://hpctoolkit.org/ http://hpctoolkit.org/ http://hpctoolkit.org/ http://www.mpcci.de/multiphysics-engineering.html http://www.npcci.de/multiphysics-engineering.html http://www.tddft.org/programs/octopus/wiki/index.php/Main_Page http://www.tddft.org/programs/octopus/wiki/index.php/Main_Page
379 380 381 383 383 384 385 388 389 390 390 391 392 393	Rhadoop cfour chombo dakota globus globus gli hpctoolkit mongodb mpb mpcci ncl octopus pbzip pcp	performance assessment easy while providing accurate diagnoses of core, chip, and node-level performance bottlenecks. Because of this focus, it suffices (and is recommended) to use scaled-down data sets and resource configurations to obtain accurate analyses. RHadoop is a part of Hadoop, an open source data processing framework. CFOUR (Coupled-Cluster techniques for Computational Chemistry) is a program package for performing high-level quantum mechanical calculations on atoms and molecules. Chombo provides a set of tools for implementing finite difference methods for the solution of partial differential equations on block-structured adaptively refined rectangular grids. Both elliptic and time-dependent modules are included. Chombo supports calculations in complex geometries with both embedded boundaries and mapped grids, and Chombo also supports particle methods. Most parallel platforms are supported, and cross-platform self- describing file formats are included. Dakota — A Mutilevel Parallel Object-Oriented Framework for design optimization Globus provides the GSIM/Proxy in addition to GridFTP cli tools GT-SUITE offers the only true virtual engine/powertrain tool, capable of integrated simulations of the total engine and powertrain system HPCTookit is an integrated suite of tools for measurement and analysis of program performance on computers ranging from multicore desktop systems to the nation's largest supercomputers. By using statistical sampling of timers and hardware performance counters, HPCTookit collects accurate measurements of a program's work, resource consumption, and inefficiency and attributes them to the full calling context in which they occur. 7 This program computes definite-frequency eigenstates (harmonic modes) of Maxwell's equations in periodic dielectric structures for arbitrary wavevectors, using fully-vectorial and three-demensional methods. It is especially designed for the study of photonic crystals (a.k.a. photonic band-gap materials), but is also applicable to man	http://www.tacc.utexas.edu/perfexpert/ https://github.com/RevolutionAnalytics/RHadoop/wiki http://www.cfour.de/ http://commons.lbl.gov/display/chombo/Chombo+-+Software+for+Adaptive+Solutions+of+Partial+Differential+Equations http://dakota.sandia.gov/ http://dakota.sandia.gov/ http://www.globus.org/ http://hpctoolkit.org/ http://hpctoolkit.org/ http://hpctoolkit.org/ http://hpctoolkit.org/ http://www.npcci.de/multiphysics-engineering.html http://www.ncl.ucar.edu/ http://www.ncl.ucar.edu/ http://www.tddft.org/programs/octopus/wiki/index.php/Main_Page http://oppression.ca/ http://oppression.ca/

	A	S	T
		Teragrid/Xsede Client Toolkit for connecting with your Teragrid/Xsede portal password to Xsede/Teragrid resources using gsissh myproxy-login and globus-	
395	tct	url-copy.	https://www.xsede.org/user-portal
		Maximally-localised Wannier functions (MLWFs) and Wannier90, the computer program that calculates them. Both wannier90.x (the program) and	
396	wannier90	libwannier.a (the library) are built and available.	http://www.wannier.org/
		The Weather Research and Forecasting (WRF) Model is a next-generation mesoscale numerical weather prediction system designed to serve both	
		operational forecasting and atmospheric research needs. It features multiple dynamical cores, a 3-dimensional variational (3DVAR) data assimilation	
397	wrf	system, and a software architecture allowing for computational parallelism and system extensibility.	http://www.wrf-model.org/index.php
398	xsede	Xsede Client Toolkit for connecting with your Xsede portal password to Xsede resources using gsissh myproxy-login and globus-url-copy.	https://www.xsede.org/user-portal
399	cfx	ANSYS CFX is a high-performance, general purpose CFD program that has been applied to solve wide-ranging fluid flow problems.	http://www.ansys.com/Products/Simulation+Technology/Fluid+Dynamics/Fluid+Dynamics+Products/ANSYS+CFX
400	chemkin	Loads the CHEMKIN runtime environment CHEMKIN software enables the simulation of complex chemical reactions	http://www.reactiondesign.com/
401	cubit	loads the CUBIT Mesh Generation Kit	http://cubit.sandia.gov/index.html
		Dytran is an explicit finite element analysis (FEA) solution for analyzing complex nonlinear behavior involving permanent deformation of material	
402	dytran	properties, or the interaction of fluids and structures.	http://www.mscsoftware.com/
403	marc	MARC is a non-linear FEA program	http://www.mscsoftware.com/Products/CAE-Tools/Marc.aspx
404	stata-mp	Stata is a complete, integrated statistical package that provides everything you need for data analysis, data management, and graphics.	http://www.stata.com/
		Computational fluid dynamics software for complex geometries with moving boundaries. Licensing: This software is licensed for academic, noncommercial	
405	converge	use at the University of Michigan	http://convergecfd.com/
406	helios	HELIOS Generalized Geometry Lattice Analysis package. Includes Casmo and Simulate from the Studsvik suite.	http://studsvikscandpower.com/products/helios.php
		Kintecus is chemical modeling software for simulation of combustion, nuclear, biological, enzyme, atmospheric and many other chemical kinetic and	
407	kintecus	equilibrium processes. Kintecus is licensed only for use by specific classes. If you are not in one of those classes, you will not be able to run it.	
408	lahey	loads the Fujitsu Lahey LF95 environment	http://www.lahey.com/
409	mcnp5	loads the mcnp5 environment	https://mcnpx.lanl.gov/
410	molpro	Molpro is a complete system of ab initio programs for molecular electronic structure calculations	http://www.molpro.net/
411	synopsys	Synopsys Synthesis environment	http://www.synopsys.com/home.aspx