PRIYA VASHISHTA

Collaboratory for Advanced Computing and Simulations
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CURRENT POSITION: Director, DOE Materials Genome Center: *Materials Genome Innovation for Computational Software (**MAGICS**)*
Founding Director, Collaboratory for Advanced Computing and Simulations
University of Southern California
Professor of Physics & Astronomy, and Computer Science

EDUCATION: Ph.D., Indian Institute of Technology, Kanpur, India, 1967

PERSONAL: Born August 24, 1944; Wife: Annie-Claude; US Citizen; Naturalization Date: May 1, 1995

PROFESSIONAL EXPERIENCE

2015 - Present Director, DOE Materials Genome Center: *Materials Genome Innovation for Computational Software (**MAGICS**)*
2002 - Present Director, Collaboratory for Advanced Computing and Simulations (CACS), and Professor of Chemical Engineering & Materials Science, Physics and Astronomy, and Computer Science, University of Southern California
1990 - 2002 Founding Director, Concurrent Computing Laboratory for Materials Simulations (CCLMS), Louisiana State University, Baton Rouge
1990 - 2002 Cray Professor of Computational Sciences, Louisiana State University, Baton Rouge
1984 - 1990 Senior Scientist, Argonne National Laboratory
1979 - 1982 Director, Solid State Science Division, Argonne National Laboratory
1972 - 1979 Member of the Research Staff, Argonne National Laboratory
1971 - 1972 Assistant Professor, Western Michigan University
1970 - 1971 Assistant Professor, Northwestern University

CONCURRENT POSITIONS

1986 - 1987 Visiting Professor, Cornell Theory Center and LASSP, Cornell University with ***Kenneth G. Wilson*** (*Nobel Prize, 1982*)
1983 Visiting Professor, Institute for Theoretical Physics, UC-Santa Barbara
1982 - 1985 Adjunct Professor of Physics, Northwestern University
1976 - 1977 Visiting Professor, University of California at San Diego with ***Walter Kohn*** (*Nobel Prize, 1998*)
1976 Visiting Scientist, Bell Laboratories, Murray Hill, NJ
1972 Visiting Scientist, IBM Research Center, Yorktown Heights
1968 - 1970 Postdoctoral Research Associate, McMaster University, Hamilton, Ontario, Canada
1966 - 1968 Postdoctoral Fellow, St Andrews University, St. Andrews, Scotland

Awards and Honors

2020 The Best Paper Award, ACM International Conference on High Performance Computing in Asia-Pacific Region
2013 The Best Paper Award, IEEE International Workshop on Parallel and Distributed Scientific and Engineering Computing

2010	Associates Award for Creativity in Research, University of Southern California
2008	Engineering Faculty Senior Research Award, University of Southern California
2002	Best Paper, IEEE Virtual Reality Conference
2001	The Best Technical Paper Award, IEEE/ACM Supercomputing Conference
1999	Fellow of the American Physical Society
1997	Sustained Excellence Performer Award in Ultra Dense, Ultra Fast Computing Components, DARPA
1997	Sustained Excellence Performer Award in Ultra Dense, Ultra Fast Computing Components, DARPA
1990	United Nations Development Program Fellowship Award for Research and Consultation in India
1989	Japan Society for the Promotion of Science Senior Fellowship Award for Research in Japan,
1985	Brazilian Science Research Council Fellowship Award for Research in Brazil
1985	Japan Society for the Promotion of Science Senior Fellowship Award for Research in Japan
1976	The University of Chicago Award for Distinguished Performance at Argonne National Laboratory

TOTAL PUBLICATIONS: 500+

Edited books: 11

Book chapters: 13

INVITED TALKS: 350+

VISION:

Computing technology will grow by a factor of more than a thousand in the next ten years. Our goal is to follow this computing revolution from petaflops (10^{15} flops) to exaflops (10^{18} flops). Using this unprecedented computing power, available for the first time in the history of science and engineering, it will be possible to carry out realistic simulations of complex systems and processes in the areas of materials genome, nanostructured materials and processes, nanotechnology, and bioengineered systems. Coupled with AI and machine learning, this will offer unprecedented opportunity for research as well as modifying graduate and undergraduate education that incorporates simulations and machine learning in materials science & engineering, and physics.

COLLABORATORY FOR ADVANCED COMPUTING AND SIMULATIONS (CACS)

The CACS has two main objectives: i) multidisciplinary research involving computational grand challenges in materials; and ii) the development of undergraduate and graduate education in computational science and engineering.

RESEARCH INTERESTS

Our primary research interests are in multiscale materials simulations that combine molecular dynamics with density functional theory for reactive processes and finite element for continuum behaviors on petascale supercomputers and as well as on future exascale computers. We are also designing multi-resolution algorithms and machine learning tools for large-scale quantum dynamics simulations. Current simulation effort focuses on: 1) Nano-bubble collapse in water under shock compression, shock propagation in ceramics and self-assembled monolayers; 2) reactive and quantum dynamics of hydrogen

production from Al-Li nano-particles in water; 3) nano-engineered energetic materials and munitions, dynamic friction under extreme conditions; 4) nucleation and growth of cracks and fracture, stress corrosion, and delamination at interfaces; 5) self-organized monolayers (SAM) and hybrid physical and biological systems consisting of nanostructures linked with membranes, DNA and peptides.

EDUCATIONAL PROGRAMS

Dual-degree Program

We have established a unique ***dual-degree program***, in which students can obtain a Ph.D. in the physical sciences or engineering and an M.S. in computer science with emphasis on data science in five years. The program is designed to produce a new generation of computational scientists and engineers who will make innovative use of emerging information technologies to address Grand Challenge problems in their application domains. We have introduced cross-disciplinary courses in support of this dual-degree program. MS program in computer science accessible to students anywhere and anytime, exploiting the infrastructure of USC's top-rated distance-education learning program.

Courses Taught

Basics of atomistic simulation of materials;

Molecular dynamics simulations of materials and processes;

Thermodynamics and statistical physics;

Solid-state physics; freshmen academy.

OUTREACH

We have established a special program for undergraduate students and their faculty mentors from Historically Black Colleges and Universities (HBCUs) and Minority Serving Institutions (MSIs). Groups of two students and one faculty are brought together from 12 institutions to participate in a one-week intense hands-on workshop on computational sciences and engineering. Students come with varying backgrounds—freshmen through seniors—and receive hands-on experience in parallel computing, including the assembly of Linux nodes from off-the-shelf components, loading them with scientific and simulation software, and connecting them to a Gigabit switch. This Linux parallel cluster is then used for algorithmic and simulation exercises in a tutorial setting. In addition to hands-on activities, the participants visit research laboratories and attend lectures on emerging trends in physical and computer sciences and engineering. Follow-on activities with participants include: (1) Bringing students back for summer research; (2) remote research experiences for students facilitated by the loan of the Linux nodes they build at the workshop; and (3) inviting 2-3 students and 2-3 faculty mentors from the previous workshop to help run the next workshop.

The latest workshop for underrepresented groups "***3rd MAGICS Workshop on Materials Software***", was held on Nov. 11 – 13, 2018 at the Gaithersburg Marriott Washingtonian Center, Gaithersburg, Maryland, It was jointly organized with Howard University, chartered by U.S. Congress on March 2, 1867, shortly after the end of the civil war, is a historically black university in Washington, D.C.

RESEARCH INFRASTRUCTURE

***Local parallel computing resources*:** We established a Collaboratory for Advanced Computing and Simulations (CACS) at USC in September 2002. The CACS has a 4,192-processors.

***Collaborative immersive and interactive visualization facilities*:** The CACS also has: 1) a 14'×8' tiled display driven by a Linux cluster and 2) an immersive and interactive virtual environment, ImmersaDesk. The ImmersaDesk provides interactive, stereoscopic data projection. The Viz Lab consists of a dedicated

space with immersive audio and visualization capabilities for tutorials, lectures, meetings, conferences, and all other forms of interactions that currently require face-to-face meetings.

***National and international computing resources*:** We have access to the IBM Blue Gene/Q and Intel Theta at Argonne National Labpratory via one of the largest DOE-INCITE grants for our research projects.

CONCURRENT COMPUTING LABORATORY FOR MATERIALS SIMULATION (*CCLMS*) AT LSU

We established a *Concurrent Computing Laboratory for Materials Simulations (CCLMS)* in 1990 at LSU i) to conduct multidisciplinary research in computational materials sciences and ii) to develop interdisciplinary educational programs involving high performance computing and communications.

CCLMS Computational facilities: With $3 million in infrastructure enhancement grants from the State of Louisiana and federal agencies, the *CCLMS* is equipped with a number of parallel architectures—a Linux cluster connected by a hierarchical network fabric consisting of high speed switches; a Digital Alpha cluster on Gigaswitches; and a 64-cell Intel iWarp systolic array architecture. *We brought the parallel clusters and Intel iWarp to USC in 2002.*

Visualization facilities: An immersive and interactive 3D virtual environment called *ImmersaDesk*; a multiprocessor SGI Onyx2 with an InfinityReality2 graphics pipeline; an SGI Octane MXE; and a number of SGI graphics workstations. *We also brought the ImmersaDesk to USC.*

Biological Computation and Visualization Center: In 2000 we received a five-year, $3.9 million competitive grant from the State of Louisiana to establish a Biological Computation and Visualization Center (BCVC). It involves faculty, postdocs, and graduate students from the biological, physical, and computer sciences, and engineering at Louisiana State University and A&M College in Baton Rouge and Louisiana State University Health Sciences Center in New Orleans. Multidisciplinary research activities within the *BCVC* focus on massively parallel simulations, data mining, and immersive and interactive visualization of biological applications. An important educational goal of the *BCVC* is to propel students into careers in emerging areas of biological and information technologies both in academic and industrial settings. In 2002, the BCVC also received a three-year, $5.9 million grant from the National Institute of Health.

ADMINISTRATIVE EXPERIENCE

From 1979-82, I was the Director of the Solid State Science Division at Argonne National Laboratory. The division consisted of scientific and technical personnel, including 50 Ph.D.-level Principal Investigators. My job involved securing block funding from the Department of Energy and overseeing various research and instrument development programs. During my tenure, the division was also involved in the establishment of a DOE supported $50 million facility called the Intense Pulsed Neutron Source (IPNS). The IPNS, during its operation, was recognized as one of the best pulsed-neutron research facilities in the world. The IPNS facility is now closed after a 30+ years of very successful operation.

MULTIINSTITUTIONAL, MULTIDISCIPLINARY ACTIVITIES

Completed_NSF-ITR (Information Technology Research), "De Novo Hierarchical Simulations of Stress Corrosion Cracking in Materials": We are spearheading (Kalia, Nakano, Vashishta—PI) a $3.8M project with Caltech (W. Goddard and M. Ortiz) and Purdue (A. Grama) to study atomistic mechanisms of stress corrosion cracking by high-end parallel and Grid simulations that seamlessly integrate quantum mechanical simulations, molecular dynamics simulations with reactive and nonreactive force fields, accelerated dynamics, and atomistically informed continuum models.

Completed_NSF-PetaApps, "Petascale Hierarchical Simulations of Biopolymer Translocation through Silicon Nitride and Silica Nanopores and Nanofluidic Channels": We are spearheading (Kalia, Nakano, Vashishta—PI) a $2M project with UC Santa Barbara (H. Metiu), Utah (M. Hall) and Purdue (A. Grama) to perform hybrid quantum mechanical-atomistic-mesoscale simulations of ion transport and translocation of biopolymers such as DNA and RNA through nanometer scale pores and channels in silica and silicon nitride membranes.

Completed_NSF-EMT (Emerging Models and Technologies for Computation), "Petascale Simulations of DNA Dynamics and Self-Assembly": We are spearheading (Kalia, Nakano, Vashishta—PI) this EMT project with Purdue (A. Grama) to develop a predictive hierarchical petascale simulation framework to study DNAs through solid-state nanopores and shear-induced DNA self-assembly.

Completed_DOE-SciDAC, "Hierarchical Petascale Simulation Framework for Stress Corrosion Cracking": We (Kalia, Nakano, Vashishta—PI) are leading this collaboration with Kaxiras (Harvard) G. Lu (Cal State); A. Grama (Purdue); J. Moriarty and L. Yang (Livermore); A. Voter (Los Alamos). The project focuses on scalable parallel and distributed computational framework for stress corrosion cracking.

Completed_DOE SciDAC-e, "*Performance enhancement of simulating the dynamics of photoexcitation for solar energy conversion*": In collaboration with Maryland (J. Hollingsworth), LLNL (D. Quinlan), and Utah (M. Hall), we (R. Lucas—PI, J. Chame, P. Diniz, R. K. Kalia, A. Nakano, P. Vashishta) enhance the performance of our petascale nonadiabatic quantum molecular dynamics simulation codes to study photoexcitation dynamics for solar energy applications.

Completed_DOE-EFRC (Energy Frontier Research Center) "Emerging Materials for Solar Energy Conversion and Solid State Lighting": This $12M EFRC involves collaboration between University of Southern California, University of Illinois, University of Michigan and University of Virginia. Research at the Center is focused on emerging materials for solar energy and solid-state lighting and the invention of new solar cell and LED designs based on nanostructured and organic materials. Multimillion atom simulations of nanostructured materials will be carried out by Kalia, Nakano and Vashishta at USC.

Completed*_NSF—Cholesterol *Flip-flop Dynamics and Nanomechanical Response of Deformed Biomembranes: Experiments and Petascale Simulations: In this joint experimental and computer simulation project, we (Kalia, Malmstadt, Nakano, Vashishta at USC) are investigating the nanomechanical response of phase separated liquid-ordered and liquid-disordered domains to membrane deformation and how cholesterol flip-flop dynamics between the membrane leaflets affect the mechanical response.

Completed_NSF CDI-Type II: "Probing Complex Dynamics of Small Interfering RNA (siRNA) Transfection by Petascale Simulations and Network Analysis": We (Kalia, Nakano, Vashishta at USC; Grama at Purdue) are performing multimillion-atom simulations to study: (1) the effect of siRNAs on the molecular structure of lipid membranes and how structural changes affect the membrane permeability; and (2) delivery of siRNAs encapsulated in liposomes by ultrasound.

***Completed*_DoD-ARL**: "Advanced modeling and characterization of strategic materials," with Temple
(September 29, 2016 to September 28, 2018) P. Vashishta, (Principal), R. K. Kalia, A. Nakano
USC portion $1,000,000

***Completed*_DTRA:** "Molecular Mechanisms of Spore Killing by Corrosive and Detonation Product Gases: Reactive Molecular Dynamics Coupled with Graph-Theoretic Methods"
Priya Vashishta, Rajiv Kalia, Aiichiro Nakano (USC)
DoD-Defense Threat Reduction Agency: $750,000 for 5 years

***Completed*_AFOSR:** "Real-time dynamics of hot spots in microstructured energetic materials: experiments and simulations"
Dana Dlott and Kenneth Suslick (UIUC), Priya Vashishta, Rajiv Kalia, Aiichiro Nakano (USC)
DoD-Air Force office of Scientific Research: $750,000 for 5 years

***Completed*_NSF:** "Probing complex dynamics of small interfering RNA (siRNA) transfection by petascale simulations"
P. Vashishta, R. K. Kalia, A. Nakano (USC)
National Science Foundation, Collaborative Research: CDI-Type II, $1,080,000 for 4 years

***Completed*_ONR:** "Synthesis, characterization, and multiscale simulations of low-cost high density insensitive high performance energetic composite molecular systems"
P. Vashishta, R. K. Kalia, A. Nakano, K Christe (USC); D. S. Stewart (UIUC); Y. Gupta (WSU)
DOD_Office of Naval Research, $3,175,000 for 5-years

CURRENT RESEARCH FUNDING

1. **DOE-BES** Materials Genome Center at USC: "Computational Synthesis of Materials Software Project with Validation on Layered Low Dimensional Functional Materials and Ultra Fast X-Ray Laser Experiments"
Priya Vashishta-PI, Malancha Gupta, Rajiv K. Kalia, Aiichiro Nakano, Oleg Prezhdo (USC), P. Ajayan, (*Rice),* Uwe Bergmann and David Fritz (Stanford), William A. Goddard, III (Caltech), Kristin A. Persson (LBL), David J. Singh, (*ORNL)*
DOE-Basic Energy Sciences: $8,000,000 for 4 years

2. **DOE-Materials Science**: "Self-Healing Nanomaterials: Multimillion-Atom Reactive Molecular Dynamics Simulations"
P. Vashishta—PI, R. Kalia, A. Nakano (USC)
DOE-Materials Science & Engineering: $540,000 for 3 years

3. **ONR:** "Synthesis, characterization, and multiscale simulations of low-cost high density insensitive high performance energetic composite molecular systems"
P. Vashishta, R. K. Kalia, A. Nakano
DOD_Office of Naval Research, $1,000,000 for 5-years

PENDING

4. DOE: "Accelerating Diffuse Neutron Scattering Research on Far-from-Equilibrium Quantum Phenomena Using Exascale Neural Network Quantum Molecular Dynamics (Ex-NNQMD)"

The overarching goal of this SciDAC Center is to develop a transformative simulation-AI-diffuse neutron scattering (DNS) approach to determine and control far-from-equilibrium topological structures and their excitations in a new class of "polar quantum materials".
DOE_BES: $8,000,000.00

5. NSF: "Quantum Materials Dynamics at the Nexus of Quantum Computing, Artificial Intelligence and Exascale Computing"
The goal is to develop software and training infrastructures to facilitate the adoption of quantum computational science (QCS) by a broad community for solving challenging scientific problems using quantum computers.
NSF-CISE_$450,000.

6. NSF: "CyberMAGICS - Cyber Training on Materials Genome Innovation for Computational Software"
The goal of this project is to train a new generation of materials cyberworkforce, who will solve challenging materials genome problems through innovative use of advanced cyberinfrastructure (CI) at the exa-AI-QC nexus.

NSF-CISE_$450,000.

7. NSF: "New Computational and Experimental Approaches to Assess the Fundamental Mechanisms of Creep"
This proposal will investigate the fundamental basis for the phenomenological equations for creep in crystalline materials with focus on high temperature plasticity.
NSF-ENG_$450,000.

8. NSF: "Chalcogenide quantum materials: From concept to scalable cyber manufacturing through AI-guided processing"
The role of materials R&D in the economy is enormous but not widely recognized. Pervasive technologies like efficient internal combustion engines rely on the results of materials engineering (e.g. sensors, nanostructured alloys) and depend on materials manufacturing.
NSF-DMR_$560,000.

8. NSF: "New Computational and Experimental Approaches to Assess the Fundamental Mechanisms of Creep and Embrittlement in Additively Manufactured IN625 and IN718"
The objective of this proposal is to evaluate the suitability of additively manufactured (AM) superalloys to replace such alloys as cast and wrought IN625 and IN 718.
ONR_$450,000.

SERVICE AT USC

High Performance Computing and Communications (HPCC) at USC

1. Faculty Advisory Council on High Performance Computing and Communications Committee, Chair, 2004-present
2. Cyberinfrastructure Architecture Committee, 2004-2005

USC Wide Committees:

3. Provosts Committee on ISD Reorganization, 2005-2006
4. Joint USC Senate and Provost Committee on Research, 2005-2006
5. Provost Strategic Planning Committee, 2008-2009
6. Graduate School Committee on Fellowships, Prizes, and Awards, 2008-2010
5. Provost Strategic Planning Committee, 2009-2011
7. Applied Mathematics task force, 2009

School of Engineering:

8. Provost Committee for the Search of Dean of School of Engineering, 2005-2006
9. Engineering Faculty Council, 2004-2006
10. Committee for the Merger of Chemical Engineering and Materials Science, 2004-2005
11. Green Engineering Curriculum Committee, 2010-2013
12. APT Committee, 2010-2015
13. Transformative Faculty Hire Committee, Chair, 2010-2015

Mork Family Department of Chemical Engineering and Materials Science

14. Chemical Engineering Centennial Committee, 2005-2006
15. EFC appointed MFD Chair evaluation committee, 2009
16. Merit review Committee, 2008-2015
17. Ping Lecture committee, 2009

18. Spitzer Lecture committee, Chair, 2012, 2013, 2014
19. MFD Executive Committee, 2011-2014

Dornsife College of Letters, Arts and Sciences :

20. College Initiatives Advisory Committee, 2003-2004
21. Physics and Astronomy Executive Committee, 2005-2006

CONFERENCE ACTIVITIES [CONFERENCE CHAIRMAN/CO-CHAIRMAN]:

- 26th "Midwest Solid State Conference", Argonne National Laboratory, November 1978
- International Conference on "Fast Ion Transport in Solids", Lake Geneva, Wisconsin, May 1979
- International Conference on "Ordering in Two Dimensions", Lake Geneva, Wisconsin, May 1980
- 29th Midwest Solid State Conference on "Novel Materials and Techniques in Condensed Matter", September 1981
- 9th Midwest Solid State Theory Symposium on "Melting, Localization and Chaos", November 1981
- Workshop on "Phase Transitions and Gauge Theories," Argonne National Laboratory, July 19 - September 10, 1982
- "Rahman Festschrift", Argonne National Laboratory, November 12-13, 1984
- X Pan American Workshop on "Condensed Matter Theories," Argonne National Laboratory, July 21-26, 1986
- "Highlights in Condensed Matter Physics", International Center for Theoretical Physics, Trieste, Italy, August 1-3, 1989
- "Concurrent Computing in the 90's", Louisiana State University, August 17, 1990
- "Undergraduate and Graduate Education in Computational Sciences", Louisiana State University, April 29-30, 1991
- "Parallel Computing Adventure: An Enrichment Course for High School Teachers", Louisiana State University, August 3-7, 1992
- International Conference on "Concurrent Computing in the Physical Sciences", Louisiana State University, February 18-20, 1993
- International Conference on "Toward Teraflop Computing and New Grand Challenge Applications", Louisiana State University, February 10-12, 1994
- Symposium on "Grand Challenges in Computer Simulations" at the High Performance Computing '94 of the Society of Computer Simulation in La Jolla, CA, April 11-15, 1994.
- International Conference on "High Performance Computing Technologies and Scientific Applications", Louisiana State University, February 23-25, 1995
- "High Performance Computing Symposium" at The Simulation Multiconference, The Society of Computer Simulation, Phoenix, AZ, April 9-13, 1995
- "XIX International Workshop on "Condensed Matter Theories", Caracas, Venezuela, June 12-17, 1995
- "Computational Physics Workshop for Undergraduate Students", Louisiana State University, July 17-28, 1995
- "Materials Theory, Simulation, and Parallel Algorithms", A five day Symposium at the Annual Meeting of the Materials Research Society, Boston, MA, November 27-December 1, 1995

- International Conference on " Experimental and Simulation Challenges in Nanostructured Materials", Louisiana State University, February 15-17, 1996
- International Materials Research Congress Symposium on "Computational Materials Science: Theory and Simulation of Nanostructured Materials", Cancun, Mexico, September 1-6, 1996
- International Conference on "Multiscale Phenomena in Science and Engineering", Louisiana State University, February 7-9, 1997
- International Conference on "Computer-Aided Design of High-Temperature Materials", Santa Fe, NM, July 30-August 2, 1997
- International Materials Research Congress Symposium on "Theory and Computer Simulation of Materials: Electronic Structure and Mechanical Behavior", Cancun, Mexico, September 1-5, 1997
- International Conference on "Materials and Microsystems for Extreme Environments: Experimental and Computational Challenges", Louisiana State University, February 19-21, 1998
- International Conference on "Thermo-Mechanical and Electrical Properties of High-Temperature Materials", Maui, HI, January 4-9, 1999
- Workshop on "Parallel Algorithms, Computational Efficiency and Multiscale Materials Simulations", April 2-3, 1999, New Orleans, LA
- International Materials Research Congress Symposium on "Theory and Computer Simulation of Materials - Matching of Length Scales in Materials Science: Electronic, Atomic and Continuum Properties of Materials", Cancun, Mexico, August 29-September 2, 1999
- International Conference on "Materials Design: Experimental and Computational Challenges", Louisiana State University, March 2-4, 2000
- International Conference on "Multiscale Materials Phenomena in Harsh Environments", Limassol, Cyprus, June 19-24, 2000
- International Conference on "Multiscale Simulation, Theoretical, and Experimental Approaches to Deformation, Friction, Fatigue, and Fracture", Louisiana State University, February 22-24, 2001
- "Computational Science Workshop for Underrepresented Groups", Louisiana State University, January 4-10, 2002
- Mardi Gras 2002 Conference on "Nanotechnology at the Interface of Information Technology" Louisiana State University, February 7-9, 2002
- "35th Year Anniversary of the Materials Science Department and the Inauguration of the Collaboratory for Advance Computing and Simulations at USC", University of Southern California, November 8, 2002
- "Computational Science Workshop for Underrepresented Groups", Louisiana State University, January 5-11, 2003
- Mardi Gras 2003 Conference on "Grid Computing and Simulation at the Nano-Bio Interface", Louisiana State University, February 27-March 1, 2003
- "Computational Science Workshop for Underrepresented Groups", University of Southern California, January 5-10, 2004
- "*Computational and Experimental Challenges in Physical, Chemical and Biological Systems*", August 20-21, 2004, Univ. of Southern California, Los Angeles, CA.
- "*Computational Science Workshop for Underrepresented Groups*",January 5-11, 2005, University of Southern California, Los Angeles, CA.
- "*Computational Science Workshop for Underrepresented Groups*", January 5-11, 2006, University of Southern California, Los Angeles, CA.

- "*Computational Science Workshop for Underrepresented Groups*", January 3-10, 2007, University of Southern California, Los Angeles, CA.
- "*Computational Science Workshop for Underrepresented Groups*"
  May 18-25, 2008, University of Southern California, Los Angeles, CA.
- "Scientific Discovery through Advanced Computing and Experiments", International Conference on Computing & Experimental Engineering and Sciences, April 8-13, 2009, Phuket, Thailand
- "*Computational Science Workshop for Underrepresented Groups*"
  June 21-29, 2009, University of Southern California, Los Angeles, CA.
- USC-DOE conference on "*Emerging Trends in Materials Simulations and Experiments*"
  March 24-26, 2010, Terranea Resort in Greater Los Angeles, CA.
- USC-DOE conference on "*Materials for Energy Applications - Experiment, Modeling and Simulations*"
  March 30-April 1, 2011, Terranea Resort in Greater Los Angeles, CA.
- USC-DOE conference on "Materials Genome: Simulations, Synthesis, Characterization and Manufacturing"
  April 4-6, 2012, Terranea Resort in Greater Los Angeles, CA.
- Symposium on "DNA Directed Self-assembly of Nanoparticles into Meta Materials for Energy and Other Applications"
  European Materials Research Society Spring Meeting, May 14-18, 2012, Strasbourg, France
- Western Regional Workshop on the White House Office of Science and Technology, Materials Genome Initiative
  April 7, 2014, Los Angeles, CA.
- MAGICS Center 1st Workshop: Software for Materials Science,
  University of Southern California, November 12-15, 2017
- MAGICS Center 2 nd USC Workshop Materials Software,
  University of Southern California, March 2-4, 2018

  MAGICS Center 3rd Workshop on Materials Software with the assistance from Howard University, Held at Gaithersburg Marriott Washingtonian Center, Gaithersburg, MD 20878
  November 11-13, 2018
- SPETSES Symposium on "Materials Genome Towards Exascale"
  Spetses island in Greece, June 10-15, 2018
- SPETSES Symposium on Machine Learning in Materials Genome,
  Spetses island in Greece, Greece, 15–21 June 2019
- American Physical Society Division of Computational Physics Focused Sessions "Emerging Trends in MD Simulations and Machine Learning"
  APS March Meeting, March 15–19, 2021

BOOKS, BOOK CHAPTERS, AND ARTICLES:

BOOKS:

1. Fast Ion Transport in Solids, (1979), Elsevier North-Holland, New York,
   Editors: P. Vashishta, J. N. Mundy and G. K. Shenoy
2. Novel Materials and Techniques in Condensed Matter, (1982), Elsevier North-Holland, New York
   Editors: G. W. Crabtree and P. Vashishta
3. Melting, Localization and Chaos, (1982), Elsevier North-Holland, New York
   Editors: R. K. Kalia and P. Vashishta
4. Condensed Matter Theories, Vol. 2, (1987), Plenum Press, New York
   Editors: P. Vashishta, R. K. Kalia, and R. Bishop
5. Correlations in Electronic and Atomic Fluids, (1990), World Scientific, Singapore
   Editors: P. Jena, R. K. Kalia, M. P. Tosi, and P. Vashishta
6. High Performance Computing and its Applications in the Physical Sciences, (1993), World Scientific
   Editors: D. A. Browne, J. Callaway, J. P. Draayer, R. W. Haymaker, R. K. Kalia,
   J. E. Tohline, and, P. Vashishta
7. Teraflop Computing and New Grand Challenge Applications, (1995), Nova Publishing
   Editors: R. K. Kalia and P. Vashishta
8. Condensed Matter Theories, Vol 11, (1995), Nova Publishing, New York
   Editors: E. V. Ludena, P. Vashishta, and R. F. Bishop
9. Materials Theory, Simulations, and Parallel Algorithms, (1996), Materials Research Society
   Pittsburgh, Pennsylvania
   Editors: E. Kaxiras, J. Joannopoulos, P. Vashishta, and R. K. Kalia
10. Computer-Aided Design of High-Temperature Materials, (1998), Oxford Univ. Press, Oxford
    Editors: A. Pechenik, R. K. Kalia, and P. Vashishta
11. Applying Molecular and Materials Modeling, (2002), Springer
    Editors: P. R. Westmoreland, P. A. Kollman, A.M. Chaka, P. T. Cummings,
    K. Morokuma, M. Neurock, E. B. Stechel, and P. Vashishta

BOOK CHAPTERS:

1. *Melting and Freezing in Two Dimensions: A Molecular Dynamics Study*,
   Priya Vashishta and Rajiv K. Kalia, in Melting, Localization and Chaos,
   Editors: R. K. Kalia and P. Vashishta, Elsevier North-Holland, New York (1982), p. 43.
2. *General Density Functional Theory*, Walter Kohn and Priya Vashishta,
   in Inhomogeneous Electron Gas, Editors: N. H. March (Oxford, England) and S. Lunqvist
   (Chalmers, Sweden), Plenum, New York (1983), p. 79.
3. *Molecular Dynamics Study of Superionic Conductors*, Aneesur Rahman and Priya Vashishta,
   in Physics of Superionic Conductors, Editors: J. W. Perram (Odense, Denmark), Plenum,
   New York (1983), p. 93.
4. *Theory of Electron-Hole Liquid*, Priya Vashishta, Rajiv K. Kalia and Kundan S. Singwi,
   in Electron-Hole Droplets in Semiconductors, Editors: C. D. Jeffries (Berkeley, USA),
   L. V. Keldysh (Moscow, USSR) and North-Holland, Amsterdam (1983), p. 1.

BOOKS, BOOK CHAPTERS, AND ARTICLES:

5. *New Forms of Molecular Dynamics and Superionic Conductors*
John R. Ray, Aneesur Rahman, and P. Vashishta, Superionic Solids and Solid Electrolytes--Recent Trends, Editors: A. Laskar and S. Chandra,. Academic Press, New York (1987), p.

6. *Molecular Dynamics Methods and Large-scale Simulations of Amorphous Materials*
P. Vashishta, R. K. Kalia, A. Nakano, W. Li, and I. Ebbsjö, in Amorphous Insulators and Semiconductors, Eds: M. F. Thorpe and M. I. Mitkova, Kluwer, Dordrooht (1996) p. 151-213.

7. *Structure and Mechanical Failure in Nanophase Silicon Nitride: Large-scale Molecular-Dynamics Simulations on Parallel Computers*
A. Omeltchenko, A. Nakano, K. Tsuruta, R. K. Kalia, and P. Vashishta, in Advances in Metal and Semiconductor Clusters Vol. IV: Cluster Materials, Editor: M. Duncan, JAI Press, Stamford, CN (1998), pp. 263-298.

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Structural Dynamics **8**, 014501:1-10 (2021)

508. Sulfurization of MoO_3 in chemical vapor deposition synthesis of MoS_2 enhanced by an H_2S/H_2 mixture
S. Hong, S. Tiwari, A. Krishnamoorthy, K. Nomura, C. Sheng, R. K. Kalia, A. Nakano, F. Shimojo, and P. Vashishta
Journal of Physical Chemistry Letters **12**, 1997-2003 (2021)

509. Mechanical behavior of ultralight nickel metamaterial
P. Rajak, A. Nakano, P. Vashishta and R. K. Kalia
Applied Physics Letters **118**, 081902:1-4 (2021)

510. Ex-NNQMD: extreme-scale neural network quantum molecular dynamics
P. Rajak, A. Aditya, S. Fukushima, R. K. Kalia, T. Linker, K. Liu, Y. Luo, A. Nakano, K. Nomura, K. Shimamura, F. Shimojo, and P. Vashishta
Proceedings of IPDPS Workshop on Scalable Deep Learning over Parallel and Distributed Infrastructure, ScaDL 2021 (IEEE, Portland, OR, 2021)

511. Reinforcement learning agent for chemical vapor deposition synthesis of quantum materials
P. Rajak, A. Krishnamoorthy, A. Mishra, R. K. Kalia, A. Nakano and Priya Vashishta
npj Computational Materials, accepted (2021)

512. Dielectric constant of liquid water using neural network *ab initio* molecular dynamics
A. Krishnamoorthy, K. Nomura, N. Baradwaj, K. Shimamura, P. Rakaj, A. Mishra, S. Fukushima, F. Shimojo, R. K. Kalia, A. Nakano, and P. Vashishta
Physical Review Letters **126**, 216403:1-7 (2021)

513. MISTIQS: an open-source software for performing quantum dynamics simulations on quantum computers
C. Powers, L. Bassman, T. M. Linker, K. Nomura, S. Gulania, R. K. Kalia, A. Nakano, and P. Vashishta
SoftwareX **14**, 100696: 1-6 (2021)

514. Inductive bias graph network for robust dynamics in atomistic simulation of materials
P. Rajak, A. Krishnamoorthy, R. K. Kalia, A. Nakano, P. Vashishta, and E. D. Cubuk
Proceedings of ICLR Workshop on Deep Learning for Simulations (2021)

515. Unveiling oxidation mechanism of bulk ZrS_2
L. Yang, S. C. Tiwari, S. S. Jo, S. Hong, A. Mishra, A. Krishnamoorthy, R. K. Kalia, A. Nakano, R. Jaramillo, and P. Vashishta
MRS Advances **6**, 303-306 (2021)

516. Deep reinforcement learning assisted kirigami design
P. Rajak, B. Wang, K. Nomura, Y. Luo, A. Nakano, R. K. Kalia and P. Vashishta
npj Computational Materials, accepted (2021)

517. Dielectric polymer property prediction using recurrent neural networks with optimizations
A. L. Nazarova, L. Yang, K. Liu, A. Mishra, R. K. Kalia, K. Nomura, A. Nakano, P. Vashishta, and P. Rajak
Journal of Chemical Information and Modeling **61**, 2175-2186 (2021)

518. Neural network quantum molecular dynamics to quantify and understand intermediate range order in neutron scattering experiments
P. Rajak, N. Baradwaj, K. Nomura, A. Krishnamoorthy, J. P. Rino, K. Shimamura, S. Fukushima, F. Shimojo, R. K. Kalia, A. Nakano, and P. Vashishta
Journal of Physical Chemistry Letters, accepted (2021)

INVITED TALKS AT CONFERENCES:

1. Ground State Energy Calculation of the Electron-Hole Liquid in Semiconductors
   P. Vashishta, P. Bhattacharyya and K. S. Singwi, Perspectives in Calculation of the Many-Electron Interactions in Solids, Taormina, Italy, Sept. 17-21 (1973)
2. Ground State Properties of the Electron-Hole Liquid in Ge and Si
   P. Vashishta, Symposium of the Division of Solid State Physics, American Physical Society Meeting, Philadelphia, Pennsylvania, March 25-28 (1974)
3. Surface Properties of Electron-Hole Drops in Germanium
   P. Vashishta, Oji Seminar on Physics of Highly Excited States in Solids, Tomakomai City, Hokkaido, Japan, Sept. 10-13 (1975)
4. Surface Properties of Electron-Hole Liquid in Germanium
   P. Vashishta, Midwest Theoretical Solid State Physics Symposium, Purdue University, West Lafayette, Indiana, Nov. 10-11 (1975)
5. Electron-Hole Drops in Semiconductors
   P. Vashishta, NATO Advanced Institute on Strongly Coupled Plasmas, Universite d'Orleans, Orleans-la-Source, France, July 6-23, (1977)
6. Electron-Phonon Interaction in Transition Metals
   P. Vashishta, CECAM Workshop on Electron-Phonon Interactions, Orsay, France, July 18 - Sept. 2, (1977)
7. Electron-Phonon Interaction in Transition Metals and Alloys
   P. Vashishta, Midwest Theoretical Solid State Physics Symposium, Argonne National Laboratory, Argonne, Illinois, Oct. 17-18 (1977)
8. Nature of Ionic Motion in AgI and CuI
   P. Vashishta, International Conference on Fast Ion Transport in Solids, Lake Geneva, Wisconsin, May 21-25 (1979)
9. Molecular Dynamics Studies of Superionic Conductors
   P. Vashishta, NATO Advanced Study Institute on Physics of Superionic Conductors, Odense, Denmark, Aug. 11-23 (1980)
10. Nature of Ionic Motion in Superionic Conductors
    P. Vashishta, International Meeting on Solid Electrolytes-Solids State Ionics and Galvanic Cells,Tokyo, Japan, Sept. 15-19 (1980)
11. Nature of Order-Disorder Transition in Superionic Conductors
    P. Vashishta, Symposium of the Division of Condensed Matter Physics, American Physical Society Meeting, Phoenix, Arizona, March 16-20 (1981)
12. Molecular Dynamics Study of 2D Melting
    P. Vashishta, NATO Advance Study Institute on Nonlinear Phenomena at Phase Transitions and Instabilities, Geilo, Norway, March 29-April 9 (1981)
13. Melting and Nucleation of a Two-Dimensional Coulomb Solid
    R. K. Kalia and P. Vashishta, International Conference on Physics of Intercalation Compounds, I.C.T.P. Trieste, Italy, July 6-10 (1981)
14. Melting and Freezing in Two Dimensions
    P. Vashishta, 9th Midwest Solid State Theory Symposium on Melting, Localization and Chaos, Argonne National Laboratory, Argonne, Illinois, Nov. 2-3 (1981)
15. Strategic Planning for Solid State Science Research and Free Electron Laser Facility
    P. Vashishta, 83rd Meeting of Argonne University Association Board of Trustees,

INVITED TALKS AT CONFERENCES:

Carnegie-Mellon University, July 12-13 (1982)

16. Melting, Freezing and Order-disorder Transition in 2-Dimensions
P. Vashishta, VIth Pan American Workshop on Condensed Matter Theories, Washington University, St. Louis, MO, Sept. 20-Oct. 1 (1982)

17. Mass and Charge Transport in Solid Electrodes and Electrolytes
P. Vashishta, American Institute of Chemical Engineers Annual Meeting, Los Angeles, California, Nov. 14-18 (1982)

18. Melting on Corrugated Surfaces
P. Vashishta, NATO Advanced Study Institute on Multicritical Phenomena, Geilo, Norway, April 10-21 (1983)

19. Melting and Freezing Transitions on Corrugated Surfaces
P. Vashishta, Third International Conference on Recent Progress in Many-Body Theories, Altenberg, West-Germany, August 29-September 3 (1983)

20. Anisotropic Coulomb Liquids and Structural Phase Transitions
P. Vashishta, Eighth Workshop on Condensed Matter Theories, Granada, Spain, Sept. 17-28, 1984

21. Ionic Transport and Structural Phase Transitions in Superionic Conductors
P. Vashishta, American chemical Society Symposium on "Transport Phenomena in Solids and Related Systems", Miami, April 29 - May 3, 1985

22. Glasses in Two dimensions
P. Vashishta, Annual Brazilian Physical Society Meeting, Sao Laurenco, M. G., Brazil, May 29-June 1, 1985

23. Fractal Dimensionalities of Brownian Trajectories and Brown Isosets in Two Dimensions
P. Vashishta, Workshop on Many-Body Theories, San Francisco, CA, August 5-10, 1985

24. Structure and Dynamics of Superionic Conductors
P. Vashishta, International Conference on Many-Body Theories, San Francisco, CA, August 12-17, 1985

25. Use of Computer Simulation Techniques to Study Atomic Migration in Solids
P. Vashishta, International Conf. on Solid State Ionics, Lake Tahoe, CA, August 14-24, 1985

26. Electron Glass in Two Dimensions
P. Vashishta, Annual Meeting of the Japanese Physical Society, Chiba, Japan, October 1-4, 1985

27. Structure and Dynamics of Ions in Superionic Conductors
P. Vashishta, Conference on "Strongly Interacting Condensed Matter Systems," Hakone, Japan, October 11-15, 1985

28. Molecular Dynamics Study of Low-Lying States in Glasses
P. Vashishta, Annual Meeting of the American Physical Society, Las Vegas, NV, March 31-April 4, 1986

29. Simulation of Classical and Quantum Systems
P. Vashishta, Conference on "Statistical Physics and Phase Transformations," University of Sao Paulo, Sao Carlos, Brazil, January 5-9, 1987

30. Rings and Medium Range Order in Glasses
P. Vashishta, Workshop on "Computer Simulation Studies in Condensed Matter Physics" University of Georgia, Athens, Georgia, February 8-19, 1988

INVITED TALKS AT CONFERENCES:

31. Present and Future Applications of Molecular Dynamics
    P. Vashishta, Symposium in honor of Anees Rahman, Minnesota Supercomputer Institute, University of Minnesota, Minneapolis, Minnesota, March 30-31, 1988

32. Medium Range Order in Glasses
    P. Vashishta, XII International Workshop on Condensed Matter Theories, Taxco, Mexico, August 14-20, 1988

33. Intermediate Range Order in Chalcogenide Glasses
    P. Vashishta, CECAM Planning Workshop on Computational Problems of Glasses and Networks, Amersfoort, The Netherlands, September 10-14, 1988

34. Molecular Dynamics Simulation of Mass and Charge Transport in Superionic Conductors and Structural Correlations in Chalcogenide Glasses
    P. Vashishta, Materials Society Annual Meeting, Boston, MA, Nov. 28 - Dec. 3, 1988

35. Computer Simulation of Classical and Quantum systems Using Supercomputers
    P. Vashishta, Summer School on "Condensed Matter Physics", Australian National University, Canberra, Australia, January 9-27, 1989

36. Structural Correlations and Phonon Density of States in $GeSe_2$ -- A Molecular Dynamics Study of Molten and Amorphous States
    P. Vashishta, Workshop on "Molecular Dynamics Simulations", Laguna Beach, CA, March 9-11, 1989

37. Phonon Density-of-States, Isotope Effect, and Superconductivity in $Ba_xK_{1-x}BiO_3$
    P. Vashishta, "Many Body Encounter in Minnesota", Theoretical Physics Institute, University of Minnesota, Minneapolis, MN, May 12-14, 1989

38. Molecular Dynamics Study of the Structure and Dynamics of Network Glasses
    P. Vashishta, "Special Topics in Molecular Sieve Research: Synthesis -Structure- Simulation" Argonne National Laboratory, Illinois, June 14-15, 1989

39. Nature of Phonons and Isotope Effect in $Ba_xK_{1-x}BiO_3$
    P. Vashishta, "Highlights in Condensed Matter Theory", International Center for Theoretical Physics, Trieste, Italy, August 1-3, 1989

40. Vibrational Density-of-States, Isotope Effect, and Superconductivity in $Ba_xK_{1-x}BiO_3$ Cubic Oxides
    P. Vashishta, "Recent Progress in Condensed Matter Theories", Campos do Jordao, Brazil, August 7-11, 1989

41. Nature of Phonons,Isotope Effect, and Superconductivity in $Ba_xK_{1-x}BiO_3$
    P. Vashishta, Yamada Conference on "Strongly Coupled Plasmas", Lake Yamanaka, Japan, August 29 - September 1, 1989

42. Molecular dynamics Study of the Structure and Dynamics of Network Glasses
    P. Vashishta, 7th International Conference on "Solid State Ionics", Hakone, Japan, November 5-11, 1989

43. Molecular Dynamics Simulation of Chalcogenide Glasses
    P. Vashishta, Annual American Crystallographic Association Meeting, New Orleans, April 8-13, 1990

44. Intermediate Range Order in Glasses
    P. Vashishta, Workshop on “Many Body Encounter-90”, Clemson University, Clemson,

INVITED TALKS AT CONFERENCES:

South Carolina, May 4-6, 1990

45. Molecular Dynamics Simulations of Molten and Glassy SiO_2, GeO_2, $SiSe_2$, and $GeSe_2$
P. Vashishta, International Symposium on "Molten Salts", Electrochemical Society Meeting, Montreal, May 7-11, 1990

46. Intermediate Range Order in Glasses
P. Vashishta, International Workshop on "Condensed Matter Theories", Elba, Italy, June 18-23, 1990

47. Role of Coulomb Interactions and Three-Body Forces in Network Glasses
P. Vashishta, CECAM Workshop on "Network Glasses", Orsay, France, July 2-13, 1990

48. Molecular Dynamics Study of Superionic Conductors and Network Glasses
P. Vashishta, International Conference on "Computer Applications to Material Science and Engineering", Tokyo, Japan, August 28-31, 1990

49. Computer Simulation of Glass Structures
P. Vashishta, Asian Conference on "Solid State Ionics", Beijing, China, October 29 - November 1, 1990

50. Ionic Transport in Network Glasses - A Molecular Dynamics Study
P. Vashishta, Symposium on "Solid State Ionics", Materials Research Society, Boston, November 26 - December 1, 1990

51. Nature of Phonons,Isotope Effect, and Superconductivity in $Ba_xK_{1-x}BiO_3$
P. Vashishta, Symposium on "Manifestations of the Electron-Phonon Interaction in Copper Oxides and Related Superconductors", Oaxtepec, Mexico, December 10-14, 1990

52. Crystalline Fragments in Glasses
P. Vashishta, International Symposium on "Frontiers of the Liquid and Amorphous State", Argonne National Laboratory, Illinois, August 12-15, 1991

53. Oxygen Isotope Effect in Superconducting $Ba_xK_{1-x}BiO_3$
P. Vashishta, XV International Workshop on "Condensed Matter Theories", Mar del Plata, Argentina, July 1-6, 1991

54. Computer Simulation of Network Glasses and Molecular Dynamics Algorithm on SIMD and MIMD Machines
P. Vashishta, Symposium on "Large-Scale Computations Applied to Many-Body Problems", Minnesota Supercomputer Institute, August 26-31, 1991

55. Parallel Algorithms for Molecular Dynamics Simulations on Distributed Memory MIMD Machines
P. Vashishta, Intel University Partner "Technology Focus Workshop", Timberline Lodge, Oregon, April 5-7, 1992

56. Microclusters Embedded in Zeolites: Atomistic Simulations on Parallel Architectures
P. Vashishta, "Domain-specific Methods for Massively Parallel Computational Science" Los Angeles, CA, May 14-15, 1992

57. Grand Challenges in Computer Simulation of Materials, Synthesis, and Processing
P. Vashishta, "Critical Technologies Research: Opportunities for DOE, Lawrence Berkeley Laboratory, Berkeley, CA, May 27-29, 1992

58. Molecular Dynamics Simulation of Network Glasses and MD Algorithm on Parallel (SIMD and MIMD) Architectures

INVITED TALKS AT CONFERENCES:

P. Vashishta, XVI Workshop on "Condensed Matter Theories", Puerto Rico, June 1-5, 1992

59. Structural and dynamical Correlations in Glasses
P. Vashishta, International Conference on the "Strongly Coupled Plasmas", Rochester University, August 17-21, 1992

60. Molecular Dynamics Simulations of Strongly Covalent Systems with Coulomb Interactions
P. Vashishta, International Meeting on "Computational Physics for Condensed Matter Phenomena", Tokyo, Japan, September 19-21, 1992

61. Molecular Dynamics Simulation of Network Glasses and MD Algorithm on Parallel (SIMD and MIMD) Architectures
P. Vashishta, The Second International Conference on "Computer Applications to Materials and Molecular Science and Engineering", Yokohama, Japan, September 22-25, 1992

62. Covalent Glasses at Large Positive and Negative Pressures - A Molecular Dynamics Study
P. Vashishta, "New Developments of Computer Based Design of Functional Materials", Government Industrial Research Institute, Ministry of International Trade and Industry, Osaka, Japan, September 29-30, 1992

63. Classical and Quantum Simulations of Very large Systems on Parallel Computers
P. Vashishta, Invited Talk at the Annual Meeting of Materials Research Society, Boston, November 30-December 5, 1992

64. Structure and Dynamics of Network Glasses and Crystals at Large Positive and Negative Pressures - A Molecular Dynamics Study
P. Vashishta, Invited Talk at the Annual Meeting of Materials Research Society, Boston, November 30-December 5, 1992

65. Molecular Dynamics Simulations of Covalent Glasses
P. Vashishta, Invited Talk at the American Physical Society Meeting, Oak Ridge, Tennessee, November 12-14, 1992

66. Covalent Glasses at Large Positive and Negative Pressures - A Molecular Dynamics Study
P. Vashishta, Invited Talk at the American Society for Materials, The Materials Society, Denver, Colorado, February 21-25, 1993

67. Structure and Dynamics of SiO_2 Glass at Large Positive and Negative Pressures
P. Vashishta, Invited Talk at the American Chemical Society, Denver, Colorado, March 28-April 2, 1993

68. Molecular Dynamics of Silica Under Very Large Positive and Negative Pressures on Parallel Computers
P. Vashishta, Invited Talk at the Simulation Multi Conference on High Performance Computing, Washington DC, March 29 - April 1, 1993

69. Molecular Dynamics Simulations of Classical and Quantum Systems on Parallel Computers
P. Vashishta, Computational Approaches for Novel Condensed Matter Systems, Gordon Godfrey Workshop, University of New South Wales, Sydney, Australia, July 12-17, 1993

70. Quantum Molecular Dynamics Simulations of Solid C_{60} - the Effects of Orientational Disorder and Pressure
P. Vashishta, Symposium in Honor of John Quinn's 60th Birthday, Brown University, Rhode Island, September 3, 1993

71. Molecular Dynamics and Quantum Molecular Dynamics Simulations on Parallel Architectures
P. Vashishta, 2nd International Association for Mathematics and Computer Simulation

INVITED TALKS AT CONFERENCES:

Conference on Computational Physics, St. Louis, MO., October 6-9, 1993

72. Molecular Dynamics Simulations of Structure and Dynamics of Silica at Large Positive and Negative Pressures
P. Vashishta, The Minerals Metals Materials Society Meeting, Pittsburgh, PA., Oct. 17-21, 1993

73. Materials Simulations on Parallel Architectures
P. Vashishta, Future Directions of Quantum Many Body Theory, Texas A&M University, College Station, Texas, January 14-16, 1994

74. Computer Simulation of Porous Glasses on Parallel Architectures
P. Vashishta, High Performance Computing Symposium, The Society for Computer Simulation, San Diego,CA., April 11-14, 1994

75. Vibrational Spectra of Solid C_{60} and Graphitic Tubules - A Tight Binding Molecular Dynamics Study on Parallel Computers
P. Vashishta, Electrochemical Society Meeting, San Francisco, California, May 23-27, 1994

76. Multimillion Particle Molecular Dynamics Simulation of SiO_2 and Si_3N_4 on Parallel Computers
P. Vashishta, CECAM Workshop on Computer Simulation of the Growth of Semiconducting Materials, Lyon, France, May 30-June 2, 1994

77. Classical and Quantum Molecular Dynamics Simulations on Parallel Computers
P. Vashishta, VIII International Workshop on Condensed Matter Theories, Valencia, Spain, June 6-10, 1994

78. Molecular Dynamics Simulations of Silica at Large Positive and Negative Pressures on Parallel Computers
P. Vashishta, 12th Symposium on Thermophysical Properties, Boulder, CO., June 19-24, 1994

79. Computer Simulation of Materials on Parallel Architectures - Glasses, Solid C_{60}, and Graphitic Tubules
P. Vashishta, Oji International Seminar on Elementary Processes in Dense Plasmas, Tomakomai, Japan, June 27-July 1, 1994

80. Molecular Dynamics Simulations of Covalent Amorphous Insulators on Parallel Computers
P. Vashishta, Symposium on Amorphous Insulators, Pingree Park, CO., August 7-11, 1994

81. Molecular Dynamics Simulation of High Temperature Ceramics
P. Vashishta, Workshop on Large Scale Simulations in Chemistry and Materials Science, Dayton, Ohio, September 12-13, 1994

82. Large Scale Simulation of Silica on Parallel Computers
Quartz Workshop, Naval Research Laboratory, Washington, DC, Sept. 19-20, 1994

83. Molecular Dynamics Simulations of Glasses on Parallel Computers
P. Vashishta, 1995 Simulation Multiconference, The Society of Computer Simulation, Phoenix, AZ, April 9-13, 1995

84. Million Atom Simulations of Materials on Parallel Computers - Silica, Silicon Nitride, and Ceramic Thin Films
P. Vashishta, American Ceramic Society Meeting, Cincinnati, Ohio, May 1-3, 1995

85. Studies of Nanoclusters and Amorphous Materials by Parallel Molecular Dynamics Simulations
P. Vashishta, Annual Meeting of the American Ceramic Society, Cincinnati, Ohio,

INVITED TALKS AT CONFERENCES:

May 1-3, 1995

86. Massively Parallel Computing for Materials Modeling
P. Vashishta, Fifth Conference on Computational Research on Materials, University of West Virginia, WV, May 3-5, 1995

87. Crack Propagation and Fracture in Ceramic Films -- Million Atom Molecular Dynamics Simulations on Parallel Computers
P. Vashishta, Annual Meeting of the European-Materials Research Society Meeting, Strasbourg, France, May 22-26, 1995

88. Large Scale Simulation of Amorphous Materials
P. Vashishta, XIX International Workshop on Condensed Matter Theories, Caracas, Venezuela, June 12-17, 1995

89. Simulation of Crack Propagation and Fracture in Silica and Silicon Nitride Films on Parallel Computers
P. Vashishta, Amorphous and Crystalline Insulating Thin Films III Symposium at the Annual Meeting of Solid State Devices and Materials, Osaka, Japan, August 21-24, 1995

90. Crack Propagation and Fracture Glasses and Ceramics - Million Atom Molecular Dynamics Simulations on Parallel Computers
P. Vashishta, Recent Progress in Computational Materials Science Symposium at the IVth International Conference on Advanced Materials, Cancun, Mexico, Sept. 4-8, 1995

91. Million Atom Simulations of Real Materials on Parallel Computers
P. Vashishta, SURA/EPSCoR Workshop on Computational Materials, University of Kentucky, Lexington, October 9, 1995

92. Multimillion Particle Simulations of Real Materials on Parallel Computers
P. Vashishta, Materials Modeling'95, Naval Research Laboratory, Washington, DC, October 17-18, 1995

93. Molecular Dynamics Simulations of Glasses and Ceramic Materials Using Parallel Computers
P. Vashishta, Takahashi Symposium, Solid State Ionics-10, Singapore, December 3-8, 1995

94. Amorphous and Liquid Carbon, and Dynamic fracture - Molecular Dynamics Simulations on Parallel Computers
P. Vashishta, High Performance Computing Workshop in Computational Chemistry and Materials Science, Maui High Performance Computing Center, Maui, December 21-22, 1995

95. Million Atom Simulations of Ceramic Materials on Parallel Computers - Crack Propagation, Fracture, and Sintering in Silicon Nitride
P. Vashishta, American Ceramic Society, Annual Meeting, Indianapolis, April 14-17, 1996

96. Molecular Dynamics Method and Large Scale Simulation of Amorphous Materials
P. Vashishta, NATO Advanced Study Institute, Amorphous Insulators & Semiconductors, Sozopol, Black Sea Coast, Bulgaria, May 26 - June 8, 1996

97 Million Atom Molecular Dynamics Simulation on Parallel Computers
P. Vashishta, "Theory Workshop", Advanced Photon Source, Argonne National Laboratory, Argonne, Illinois, June 24-27, 1996

98. Amorphization and Fracture in Silicon Diselenide Nanowires - A Molecular dynamics study
P. Vashishta, International Materials Research Congress, September 1-6, 1996, Cancun, Mexico

99. Large-Scale Simulation of Crack Propagation and Fracture on Parallel Computers

INVITED TALKS AT CONFERENCES:

P. Vashishta, Parallel Computing Conference, Minnesota Supercomputing Institute, Minneapolis, MN, October 3-4, 96.

100. Million Atom Simulations of Crack Propagation, Fracture, Sintering, and Densification of Ceramic Materials on Parallel Computers
P. Vashishta, Tsukuba Symposium on "Computer Simulations for Inorganic and Organic Materials and Biological Systems", Tsukuba, Japan, February 20, 1997

101. Million Atom Simulations of Crack Propagation, Fracture, Sintering, and Densification of Ceramic Materials on Parallel Computers
P. Vashishta, "International Symposium on Novel Materials", Puri, India, March 3-7, 1997.

102. Large Scale Computer Simulations of Glasses
P. Vashishta, Symposium on "Intermediate Range Order in Glasses", American Crystallographic Society of America, St. Louis, Mo, July 21-25, 1997

103. Molecular Dynamics Simulations of Si/Si3N4 Interface and Si/Si/Si3N4 Mesa on Parallel Computers
P. Vashishta, "The Second NASA Workshop on Semiconductor Device Modeling", NASA-Ames, CA, August 7-8, 1997

104. Molecular Dynamics Simulations of Si/Si3N4 Interface and Si/Si/Si3N4 Mesa on Parallel Computers
P. Vashishta, Symposium on "Theory and Computer Simulation of Materials - Electronic, Structural, and Mechanical Properties", Mexican Materials Research Society, Cancun, Mexico, September 1-4, 1997

105. Multimillion Atom Simulations of Sintering, Crack Propagation, and Fracture on Parallel Computers
P. Vashishta, "Noncrystalline Materials", Ohio State University, Columbus, Ohio, November 20, 1997

106. Multimillion Atom Simulations of Sintering, Consolidation, and Fracture of High Temperature Materials on Parallel Computers
P. Vashishta, Workshop on "Computational Materials Physics in the Southeast", Vanderbilt University, Nashville, Tennessee, November 5-8, 1997.

107. Multimillion Atom Molecular Dynamics Simulations of Si/Si/Si3N4 Nanopixel-Structural Correlations at Si/Si3N4 Interface, Fracture, and Atomic Level Stresses in the Nanopixel
P. Vashishta, Workshop on "Interfacialy Controlled Functional Materials: Electrical and Chemical Properties", Schlob Ringberg, Germany, March 8-13, 1998.

108. Large-Scale Molecular Dynamics Simulations of Mechanical Properties of Ultrafine Microstructures on Parallel Computers
P. Vashishta, Symposium on "Computational and Mathematical Models of Microstructural Evolution", Materials Research Society Meeting, San Francisco, April 13-17, 1998.

109. Multimillion Atom Molecular Dynamics Simulations of Mechanical Properties of Nanostructures in Ceramics and at Semiconductor/Ceramic Interfaces
P. Vashishta, Symposium on "Computational Modeling of Materials and Processing", American Ceramic Society, Cincinnati, Ohio, May 3-6, 1998

110 Multimillion Atom Molecular Dynamics Simulations of Ceramic Materials and Interfaces on Parallel Computers
P. Vashishta, Fourth Special Symposium on "Advanced Materials", Nagoya, Japan, May 12-14, 1998.

INVITED TALKS AT CONFERENCES:

111. Large-Scale Molecular Dynamics Simulations of Amorphous on Parallel Computers
P. Vashishta, CECAM workshop: "Intermediate Range Order in Network Forming Liquid and Glasses", Lyon, France, May, 18-20 1998.

112. Large Scale Simulations of Glasses on Parallel Computers: Nanophase Glasses and Ceramics
P. Vashishta, The 18th International Congress on Glass, San Francisco, CA, July 5-10, 1998.

113. Large Scale Computer Simulation of Glasses
P. Vashishta, Workshop on "Modeling Disordered Materials", American Crystallographic Society, Washington DC, July 18-24, 1998

114. Multimillion Atom simulations of Crack Propagation, Fracture, and Nanoindentation in Ceramic Materials
P. Vashishta, Gordon Conference on "High Temperature Materials, Processing, & Diagnostics", Plymouth, New Hampshire, July 19-24, 1998

115. Atomic Stresses in Si/Si3N4 Nanopixel - 10 Million Atom Molecular Dynamics Simulation on Parallel Computers
P. Vashishta, CECAM Workshop on "Surfaces and Interfaces far from Equilibrium", Lyon France, July 27-29, 1998

116. Atomic Stresses in Si/Si3N4 Nanopixel - 10 Million Atom Molecular Dynamics Simulation on Parallel Computers
P. Vashishta, "New Developments in High Temperature Ceramics", Istanbul, Turkey, August 12-15, 1998

117. Multimillion Atom Simulations of Silica: Crack Propagation and Fracture and Morphology of Fracture Surfaces in Amorphous Silica and Nanophase Silica
P. Vashishta, SILICA98, Mulhouse, France, September 1-4, 1998

118. Multimillion Atom Molecular Dynamics Simulations of Ceramic Materials and Interfaces on Parallel Computers
P. Vashishta, "Advanced Materials for Extreme Environments: New Experimental Opportunities in Neutron Scattering", Argonne National Laboratory, Illinois, September 11-12, 1998

119. Multimillion Atom Simulation of Crack Propagation, Fracture, and Nanoindentation in Ceramic Materials
P. Vashishta, "Computational Studies of Interfacial Phenomena: nanoscale to Mesoscale" Pacific Northwest National Laboratory, Richland, Washington, September 24-25, 1998

120. Multimillion Atom Molecular Dynamics Simulations of Ceramic Materials and Stresses in Silicon/Silicon/Silicon Nitride Nanopixel
P. Vashishta, International Union of Materials Research Society, Banglore, India, October 13-16, 1998

121. Multimillion Atom Simulations of Crack Propagation, Fracture, and Nanoindentation in Ceramic Materials and Glasses
P. Vashishta, Symposium on "Fracture and Ductile vs Brittle Behavior -Theory, Modeling, and Experiments", Materials Research Society, Boston, MA, November 30-December 4, 1998

122. Multimillion Atom Simulation of Materials on Parallel Computers -Nanopixel and Nanoindentation
P. Vashishta, "Solid State Physics Symposium 1998", Kurukshetra, India, December 27-31, 1998

INVITED TALKS AT CONFERENCES:

123. Large Scale Atomistic Simulations of Ceramic Materials and Interfaces on Parallel Computers
P. Vashishta, "Thermo-Mechanical and Electrical Properties of High-Temperature Materials", Maui,January 4-9, 1999

124. Molecular Dynamics Simulations of Materials on Parallel Computers
P. Vashishta, "Scientific Simulation Initiative", Jefferson Nat'l Laboratory, VA, January 21, 1999

125. Multimillion Atoms Simulation of High Temperature Ceramic Materials
P. Vashishta, " The Kick-off Meeting for the Eutectic Research", National Science Foundation, Arlington, VA, February 24, 1999

126. Multimillion Atom Molecular Dynamics Simulations of Ceramic Materials and Interfaces on Parallel Computers
P. Vashishta, Symposium on "Spanning the Size Scales in Materials Phenomena", American Physical Society Centennial Meeting, Atlanta, Georgia, March 22-26, 1999

127. Crack Propagation and Fracture in Nanophase Materials - Multimillion Atom Molecular Dynamics Simulation on Parallel Computers
P. Vashishta, "Nanocomposites: Design and Applications", Anchorage, Alaska, March 28-April 2, 1999

128. Multimillion Atom Simulations of Crack Propagation, Fracture, and Nanoindentation in Ceramic Materials
P. Vashishta, "High Strength Steels - Processing and Applications", Ranchi, India, April 14-15, 1999

129. Large Scale Materials Simulations on Parallel Computers
P. Vashishta, "Computational Materials Science Network on Ceramic/Polymer Interfaces", Sandia National Laboratory, Albuquerque, NM, April 22-23, 1999

130. Large Scale Simulations of Glasses on Parallel Computers: Nanophase Glasses and Ceramics
P. Vashishta, NATO Advanced Study Institute on "Physics of Glasses: Structure and Dynamics", Corsica, France, May 10-22, 1999

131. Computational Assisted Development of High Temperature Structural Materials
P. Vashishta, "DoD Challenge Workshop", Montrey, CA, June 6-10, 1999

132. Large Scale Simulations of Glasses on Parallel Computers: Nanophase Glasses and Ceramics
P. Vashishta, NATO Advanced Study Institute on "Physics of Glasses: Structure and Dynamics", Corsica, France, May 10-22, 1999

133. Multimillion Atom Molecular Dynamics Simulations of Ceramic Materials and Interfaces on Parallel Computers
P. Vashishta, "International Conference on Advanced Materials", Beijing, China, June 13-18, 1999

134. Multimillion Atom Simulations of Nanostructured Materials on Parallel Computers - Molecular Dynamics Method and Its Applications
P. Vashishta, "Computational and Applied Mathematics", St. Louis, MO, August 9-11, 1999

135. Atomistic Simulations of Nanoindentation of Silicon Nitride on Parallel Computers
P. Vashishta, "IMRC -Theory and Computer Simulation of Materials ", Cancun, Mexico, August 30-September 3, 1999

INVITED TALKS AT CONFERENCES:

136. Multimillion Atom Simulations of Nanostructured Materials on ParallelComputers - Sintering and Consolidation, Fracture, and Oxidation
P. Vashishta, "International Conference on Computational Physics ", Kanazawa, Japan, October 11-13, 1999

137. Atomistic Simulations of Nanostructures: Multimillion Atom Molecular Dynamics Simulations on ParallelComputers
P. Vashishta, "Foresight Conference on Molecular Nanotechnology ", San Jose, CA, October 15-17, 1999

138. Computational Assisted Development of High-Temperature Structural Materials
P. Vashishta, "Defense Science Board Presentation", Washington, DC, February 4, 2000

139. Multi-Million Atom Molecular Dynamics Simulations of Metal/Ceramic and Semiconductor/Ceramic Interfaces on Parallel Computers
P. Vashishta, "Workshop on Process Modeling of Laminated Multilayer Ceramic Systems", Motorola University,Tempe, Arizona, March 1, 2000

140. Molecular Simulations of Solids: Metals, Semiconductors, Ceramics, and Glasses
P. Vashishta, "International Comparative Study of Applications of Molecular and Materials Modeling", Washington, DC, March 14, 2000

141. Mulltimillion Atom Simulation of Nanostructured Materials on Parallel Computers – Sintering and Consolidation and Fracture and Oxidation
P. Vashishta, "The Materials Society (TMS)", Nashville, TN, March 12-16, 2000

142. Large-Scale Atomistic Simulations of Solid State Materials Modeling Many Millions of Atoms on Parallel Computers
P. Vashishta, "American Physical Society ", Minneapolis, March 22-24, 2000

143. Multimillion Atom Simulations of Materials on Parallel Computers - Past, Present and Future
P. Vashishta, "Celebrating the Success of LSU Computer Science" Baton Rouge, LA, April 14, 2000

144. Multimillion Atom Simulations of Nanophase Materials on Parallel Computers
P. Vashishta, "Seminars on Humanity 3000- Nanotechnology In the Future", Foundation for the Future, Bellevue, WA, April 21, 2000

145. Computer Simulations of Ceramic Interfaces
P. Vashishta, "AFOSR Meeting on Ceramic Materials and Composites" Saint Louis, MO, May 4-5, 2000

146. Multimillion Atom Simulations of Materials on Parallel Computers - Past, Present and Future
P. Vashishta, "XXIII Encontro Nacional de Física da Matéria Condensada", São Lourenço, Minas Gerais, Brasil, May 11-13, 2000

147. Multi-Million Atom Simulation of Sintering and Consolidation on Parallel Computers
P. Vashishta, "SIAM Conference on Mathematical Aspects of Materials Science", Philadelphia, PA, May 23, 2000

148. Recent Research Activities in Material Sciences at LSU: Introduction to Concurrent Computing Laboratory for Materials Simulations (CCLMS)
P. Vashishta, "Niigata University-LSU Symposium", Niigata University, Niigata, Japan, May 30, 2000

INVITED TALKS AT CONFERENCES:

149. Computational Assisted Development of High-Temperature Structural Materials
P. Vashishta, "DoD High Performance Computing Modernization Program Conference", Albuquerque, NM, June 5-8, 2000

150. Large-Scale Atomistic Simulations of Amorphous Polymers and Ceramic/Polymer Interfaces on Parallel Computers
P. Vashishta, "American CrystallographicSociety ", Minneapolis, MN, July 22-27, 2000

151. Multimillion Atom Simulation of Materials on Parallel Computers -Nanopixel, Interfacial fracture, Nanoindentation, and Oxidation
P. Vashishta, "MAPINT Symposium on Multidisciplinary Applications and Interoperable Computing", Dayton, Ohio, August 17, 2000

152. Atomistic Modeling Capabilities for Sintering of Ceramics and Fracture at Interfaces
P. Vashishta, "NSF Multi-University I/UCRC Ceramic and Composite Materials Center", Piscatway, NJ, September 27-28, 2000

153. Multimillion Atom Simulation of Nanostructured Materials – Dynamic Fracture, Nanoindentation, and Oxidation
P. Vashishta, "International Conference on Engineering & Technological Sciences 2000 – Advanced Materials", Beijing, China, October 10-14, 2000

154. Large Scale Molecular Dynamics Simulations of Materials on Parallel Computers
P. Vashishta, "ACAT2000 Conference", Fermilab, IL, October 18-19, 2000

155. Large Scale Molecular Dynamics Simulations of Nanostructured Materials on Parallel Computers– Dynamic Fracture, Nanoindentation, and Oxidation
P. Vashishta, "Joint CNRS (France)-NSF (USA) Nanomaterials Workshop – Nanomaterials Toward Engineering Applications", Montreal, Canada, October 20-25, 2000

156. Multimillion Atom Simulation of Nanostructured Materials and Devices on Parallel Computers
P. Vashishta, "DAE Solid State Physics Symposium", Guru Ghasidas University, Bilaspur, India, December 27 - 31, 2000

157. Periodic Array of Stress Domains in Silicon/Silicon Nitride Nanopixels: A Multimillion Atom Molecular Dynamics Simulation
P. Vashishta, "International Conference on Science and Technology of Nanostructured Materials", Puri, India, January 8-12, 2001

158. Multimillion Atom Simulation of Materials on Parallel Computers - Nanopixel, Interfacial Fracture, Nanoindentation, and Oxidation
P. Vashishta, "Golden Jubilee of Saha Institute of Nuclear Physics", Calcutta, India, February 5-8, 2001

159. Multimillion Atom Simulation of Nanosystems at the Forefront of Information Technology and Nanosciences, and An Integrated Dual Degree Program of Research and Education
P. Vashishta, "National Science Foundation Distinguished Lecture", Arlington, VA, February 26, 2001

160. Multimillion Atom Simulations of Mechanical Behavior of Nanostructured Materials, Interfaces, and Dynamics of Oxidation
P. Vashishta, "The Society of Materials Science, Japan (JSMS)", Osaka, Japan, May 21-26, 2001

INVITED TALKS AT CONFERENCES:

161. Hybrid Atomistic -Continuum Simulations of Nanopixels: A Multimillion Atom Molecular Dynamics Simulation
P. Vashishta, "DoD High Performance Computing", Biloxi, MS, June 18-20, 2001

162. Multimillion Atom Simulations of Materials at the Forefront of Information Technology and Nanosciences
P. Vashishta, National Research Council, National Materials Advisory Board Meeting on "Materials Research for Defense-After-Next", Woods Hole Center of the National Academy of Sciences, MA, June 27-28, 2001

163. Massively Parallel Atomistic Simulations of Nanoparticles Under Extreme Conditions
P. Vashishta, "DARPA HEDAM/NANO Workshop", Rosslyn, VA, August 6-8, 2001

164. High Temperature Materials Simulations on Parallel Computers
P. Vashishta, "Joint Annual Review of AFOSR Metallic and Ceramic Materials Program", Snowbird, UT, August 19-21, 2001

165. High Performance Computing and Visualization: Opportunities and Challenges at the Information-Bio-Nano- Interface
P. Vashishta, "Frontiers in Information Technology", Rensselaer Polytechnic Institute, Troy, NY, November 8, 2001

166. Multimillion Atom Simulations of Nanosystems on Parallel Computers
P. Vashishta, "International Workshop on Condensed Matter Theories", Canberra, Australia, December 3-8, 2001

167. High Performance Computing and Visualization: Opportunities and Challenges at the Information-Bio-Nano Interface
P. Vashishta, "Computational Sciences Workshop for Underrepresented Groups", Baton Rouge Louisiana, January 6, 2002

168. Computational Materials Science on Teraflop-to-Petaflop Computers
P. Vashishta, "DOD, DOE, NASA Joint Conference on Mission Critical Computing", Washington DC, February 4-6, 2002

169. Multimillion Atom Simulations of Nanosystems on Parallel Computers
P. Vashishta, "International Symposium on Computational Science & Engineering", Tokyo, Japan, March 5-6, 2002

170. Large-Scale Molecular Dynamics Simulations of the Oxidation of Metallic Nanoparticles
P. Vashishta, "Gordon Conference on Energetic Materials", Tilton, NH, June 19, 2002

171. Billion-atom Multiscale Simulations on a Grid
P. Vashishta, "NSF DMR Computational Materials Theory Program Review", Urbana, IL, June 20, 2002

172. Amorphous Materials in Various Incarnations
P. Vashishta, "CECAM Workshop on Atomic Structure and Transport in Glassy Networks", Lyon, France, June 24-27, 2002

173. Multimillion Atom Simulations of Nanosystems on Parallel Computers - Nanopixel, Nanoindentation and Oxidation

INVITED TALKS AT CONFERENCES:

P. Vashishta, "APS Conference on Computational Physics 2002", San Diego, August 25-28, 2002

174. Multimillion Atom Simulations of Nanosystems on Parallel Computers - Nanopixel, Nanoindentation and Oxidation of Aluminum Nanoparticles
P. Vashishta, "CNER-MSI Nanosimulation Workshop" Center for NanoEnergetics Research Minnesota Supercomputing Institute, Minneapolis, MN, August 26-27, 2002

175. Info-Bio-Nano Interface: High-Performance Computing & Visualization
P. Vashishta, "High Performance Computing 2002", Bangalore, India, December 18-21, 2002

176. Info-Bio-Nano Interface: High-Performance Computing & Visualization
P. Vashishta, "Computational Science Workshop for Underrepresented Groups", LSU@Baton Rouge, LA, January 5-11, 2003

177. Large Scale Atomistic Simulations of Reactivity of Nanosystems - Nanostructured Materials and Oxidation of Aluminum Nanoparticles
P. Vashishta, "Expanding the Envelope: Nano Materials for Aerospace Symposium" Corpus Christi, Texas, January 27-30, 2003

178. Multimillion Atom Simulation of nanosystems on Parallel Computers – nanophase Material, Nanopixel, Nanoindentation and Oxidation
P. Vashishta, "Modeling and Simulations in Micro and Nano Technologies and Materials Engineering, Toulouse, France, April 7-11, 2003

179. Multimillion Atom Simulations of Nanosystems and Interfaces on Parallel Computers
P. Vashishta, "International Conference on Non-Crystalline Inorganic Materials-Synthesis, Structure and Simulation: CONCIM", Bonn, Germany, April 8-12, 2003

180. Multimillion Atom Molecular Dynamics Simulations of Nanostructures on Parallel Computers
P. Vashishta, "International Conference on Computational Engineering & Science", Corfu, July 25, 2003

181. Multimillion Atom Molecular Dynamics Simulations of Nanostructures on Parallel Computers
P. Vashishta, "High-Tech Conference, Inha University and University of Southern California Korea" September 17-18, 2003

182. Multimillion Atom Molecular Dynamics Simulations of Nanostructured Materials
P. Vashishta, "Symposium on Multiscale Design of Nano-materials, IUMRS-ICAM2003", Yokohama, Japan, October 10-11, 2003

183. Multimillion Atom Molecular Dynamics Simulations of Nanostructures on Parallel Computers
P. Vashishta, "International Conference on Nanoscience & Technology, ICONSAT 2003", Kolkata, India, December 17-20, 2003

184. Multimillion Atom Simulations of Nanosystems and Billion Atom Walkthrough,
Priya Vashishta, "N+N Program", Washington, DC, April 15-16, 2004

185. Multimillion Atom Molecular Dynamics Simulations of Nanostructures on Parallel Computers
Priya Vashishta, "Computational Science Workshop for Underrepresented Groups", University of Southern California, 5-10January, 2004

186. Atomistic Simulations on High-end Computers and Billion Atom Walkthrough
Priya Vashishta, "Research Opportunities in Cyberengineering/Cyberinfrastructure",

INVITED TALKS AT CONFERENCES:

Drexel, April 22-23, 2004

187. Multimillion Atom Molecular Dynamics Simulations of Nanostructures on Parallel Computers
Priya Vashishta, "Atomic Scale Materials Design; Modeling and Characterization",
E-MRS Spring Meeting, Strasbourg, France, May 24-28, 2004

188. Multimillion Atom Simulations of Nanosystems and Interfaces on Parallel Computers
Priya Vashishta, Inaugural Session, 3rd International Conference "Computational Modeling & Simulation of Materials", Acireale, Sicliy, Italy, May 30-June 4, 2004

189. Billion-atom Multiscale Simulations of Nanosystems on a Grid
Priya Vashishta, NSF Division of Materials Research ITR Computational Workshop
University of Illinois, Urbana, June 17-19, 2004

190. Ultrascale Simulations High Performance Computing and Visualization
Priya Vashishta, Metaphoric Optical Computing Workshop San Diego, CA, July 12, 2004

191. Large-Scale Atomistic Simulations of Nanosystems on Parallel Computers
Priya Vashishta, "International Conference on Computational Engineering & Science",
Madeira, Portugal, July 26-29, 2004

192. Multimillion Atom Simulations and Visualization of Oxidation and Hypervelocity Impact Damage
Priya Vashishta, "Advanced Computing and Simulation", 24th Army Science Conference,
Orlando, FL, November 30, 2004

193. Multimillion Atom Simulations of Dynamics of Oxidation of an Aluminum Nanoparticle and Hypervelocity Impact Damage in AlN Ceramic,
Priya Vashishta, "Michael Klein 65th Birthday Symposium", ACS Meeting,
San Diego, CA, March 13-17, 2005

194. Multimillion Atom Simulations of Nanosystems: Structural Transformation in Nanocrystals and Hypervelocity Impact Damage
Priya Vashishta, "Algorithm and Software Developments in Nanoscale Simulations"
RISC_AIST Computational Science Workshop, Tsukuba, Japan March 22-23, 2005

195. Multimillion Atom Simulations of Hybrid Nanosystems
Priya Vashishta, "Hybrid Molecular/ Semiconductor Structures: Design of Electronics by Computation", ARO/AFOSR Workshop, Rosslyn, Virginia, June 17, 2005

196. Molecular Dynamics Simulations of High Strain Rate Deformation and Shock Propagation in Ceramics and Glasses
Priya Vashishta, "DoD-HPCMP User Group Conference" Nashville, TN, June 27-30, 2005

197. Multimillion Atom Simulations of Nanostructured Materials on Parallel Computers
Priya Vashishta, "5th International Congress of Theoretical Chemical Physics",
New Orleans, LA, July 20-26, 2005

198. Multimillion Atom Simulations of Materials Dynamics of Oxidation & Mechanical Behavior on Parallel Computers
Priya Vashishta, Workshop on "Modeling Materials in Extreme Environment (ME2)",
Washington DC, September 24-25, 2005

INVITED TALKS AT CONFERENCES:

199. Multimillion Atom Simulations of Nanosystems - Structural Transformations in Nanocrystals, Dynamics of Oxidation, and Hypervelocity Impact Damage
Priya Vashishta, Workshop on "Multiscale Modeling in Condensed Matter and Materials Science." Institute for Pure and Applied Mathematics (IPAM), UCLA, October 17-20, 2005

200. Multiscale Algorithms for Multimillion Atom Simulations on Parallel Computers
Priya Vashishta, "International Conference on Computational & Experimental Engineering & Sciences (ICCES05) ", Chennai, India, December 1-6, 2005

201. Multimillion Atom Simulations - Oxidation of Aluminum Nanoparticles, Nanophase materials, and Hypervelocity Impact Damage
Priya Vashishta, U. S. ARMY Materials Summit "Disruptive Materials Technology", Gettysburg, PA, March 14, 2006

202. Multimillion Atom Simulations of Nanostructured Materials on Parallel Computers
Priya Vashishta, "HPC Forum", Richmond, VA, April 12, 2006

203. Multimillion Atom Simulations of Reactive Nanosystems on Parallel Computers
Priya Vashishta, "11th international Ceramics Congress and 4th Forum on New Materials", Acireale, Sicily, June 4-9, 2006

204. Multimillion Atom Simulations of Dynamics of Wing Cracks and Hypervelocity Impact Damage
Priya Vashishta, " Nano-Technology and Information Technology for Space Applications", Pasadena Conference Center, Pasadena, CA, July 16, 2006

205. Multimillion Atom Simulations of Dynamics of Wing Cracks and Hypervelocity Impact Damage
Priya Vashishta, "Multiscale Nano- and Bio-Mechanics and Materials", 7th World Conference on Computational Mechanics (WCCM), Los Angeles, CA, July 16-22, 2006

206. Multimillion Atom Simulations of Dynamics of Wing Cracks and Nanoscale Damage in Glass, and Hypervelocity Impact Damage in Ceramics
Priya Vashishta, "The 12th International Symposium on Plasticity & Its Applications", Halifax, Canada, July 17-22, 2006

207. Multimillion Atom Simulations of Dynamics of Wing Cracks and Nanoscale Damage in Glass, and Hypervelocity Impact Damage in Ceramics
Priya Vashishta, "Conference on Computational Physics", Gyeongju, Republic of Korea, August 29-September 1, 2006

208. Multimillion Atom Simulations of Nanoscale Materials on Parallel Computers
Priya Vashishta, "APS California Section Meeting", CSU, Long Beach, October 27-28, 2006

209. Multimillion Atom Simulations of Nanosystems on Parallel Computers
Priya Vashishta, Chile Physical society, Santiago, Chile, November 15,-17 2006

210. Multimillion Atom Simulations of Nanosystems on Parallel Computers
Priya Vashishta, "Physics of Mesoscopic and disordered Materials", IIT Kanpur, India, December 4-8, 2006

211. Thermal Management Strategic Technology
Priya Vashishta, "Air Force Research Laboratory Workshop", Dayton, Ohio, March 29, 2007

212. Multimillion Atom Simulations of Nanostructured Materials

INVITED TALKS AT CONFERENCES:

Priya Vashishta, "5th International Symposium on Theory of Atomic and Molecular Clusters ", Richmond, Virginia, May 17, 2007

213. Multimillion to Billion Atom Simulations of Nanostructured Materials Under Extreme Conditions
Priya Vashishta, "Multi-scale and Large-scale Simulations in DoD Materials Science", Mitre Corporation, Mclean, Virginia, June 14-15, 2007

214. Multimillion to Billion Atom Simulations of Materials
Priya Vashishta, "Bold Predictions in Theoretical Chemistry: A Symposium in Honor of One of the Boldest, Bill Goddard, on the Occasion of his 70th Birthday", ACS Meeting, Boston, Mass, August 22, 2007

215. Multimillion to Billion Atom Simulations of Nanostructured Materials Under Extreme Conditions
Priya Vashishta, "Dreams, Creation and Realization of Materials Saving the Humankind", 17th IKETANI International Conference, Tokyo, Japan, September 5-8, 2007

216. Multimillion to Billion Atom Simulations of Nanostructured Materials Under Extreme Conditions
Priya Vashishta, "Ultrascale Hierarchical Simulations on High-End Parallel and Distributed Computers" , APCOM'07-EPMESC XI, Kyoto, Japan, December 03, 2007

217. Multimillion to Billion Atom Simulations of Reactive Nanosystems
Priya Vashishta, "Workshop on Pseudopotentials and Complex Materials", Jim Chelikowsky's 60th Birthday Celebration, Austin, Texas, April 10-11, 2008

218. Multimillion to Billion Atom Simulations of Nanosystems Under Extreme Conditions
Priya Vashishta, "Fragmentation Processes in the Earth" , The 21st Kronsberg Seminar, Norway, May 7-9, 2008

219. Multimillion to Billion Atom Simulations of Nanoenergetic Materials
Priya Vashishta, "Microthermal Initiator & NanoEnergetic Device Workshop", Army research Laboratory, Adelphi, MD, May 22-23, 2008

220. Multimillion Atom Simulations of Energetic Materials: Combustion of Al Nanoparticles and Reactive Nanojets in Nanostructure-Enhanced Chemical Reactions
Priya Vashishta, "Gordon Conference on Energetic Materials", Tilton, NH, June 15-20, 2008

221. Multimillion to Billion Atom Simulations of Nanostructured Materials Under Extreme Conditions
Priya Vashishta, "The 1st International Symposium on Advanced Microscopy and Theoretical Calculations", Nagoya, Japan, June 29-30, 2008

222. Multimillion to Billion Atom Simulations of Reactive Nanosystems and Stress Corrosion
Priya Vashishta, "Corrosion – Aqueous: Mechanisms, Methods and Models", Colby-Sawyer College, New London, NH, July 20-25, 2008

223. Multimillion to Billion Atom Simulations of Nanosystems Under Extreme Conditions
Priya Vashishta, "Conference on Current Trends in Computational Chemistry 2008", Jackson, MS, October 30 - November 1, 2008

224. Multimillion to Billion Atom Simulations of Nanosystems Under Extreme Conditions
Priya Vashishta, "Computational Approaches and Applications in Earth Materials Studies", American Geophysical Union, San Francisco, CA, December 15-19, 2008

225. High Performance Computing for Billion Atom Simulations of Nanosystems

INVITED TALKS AT CONFERENCES:

Priya Vashishta, "Scientific Discovery through Advanced Computing", ICCES'2009, Phuket, Thailand, April 8, 2009

226. Multimillion Atom Simulations of Reactive Nanosystems
Priya Vashishta, "Symposium on Nano-scale Energetic Materials: Fabrication, Characterization and Molecular Modeling", E-MRS 2009 Spring Meeting, Strasbourg, France, June 8-12, 2009

227. Multimillion to Billion Atom Simulations of Nanostructured Materials Under Extreme Conditions
Priya Vashishta, "Symposium in Honor of Professor Sidney Yip", USNCCM 10, Columbus, OH, July 16-18, 2009

228. Multimillion Atom Simulations of Reactive Nanosystems
Priya Vashishta, "International Workshop on Condensed Matter Theories: CMT-33", Quito, Ecuador, August 16-21, 2009

229. Fast Reaction Mechanism of Al/Al_2O_3 Nanoparticles in Oxygen: Multimillion Atom Simulations of Nanoenergetic Materials
Priya Vashishta, "Advancements in Energetic Materials & Chemical Propulsion", 8-ICICP, Cape Town, South Africa, November 2-6, 2009

230. Multimillion Atom Simulations of Nanostructured Materials
Priya Vashishta, "CIFAR Nanoelectronics Meeting", Quebec, Canada, November 16, 2009

232. Multimillion Atom Simulations of Reactive Nanosystems
Priya Vashishta, "Multiscale Modeling and Simulations of Hard and Soft Materials", JNCASR, Bangalore, India, December 17-20, 2009

233. Multimillion Atom Simulations of Reactive Nanosystems
Priya Vashishta, "Plasticity 2010", St. Kitts, January 3-8, 2010

234. Hydrogen Production from Water Using Aluminum Superatoms
Priya Vashishta, "Theory of Atomic & Molecular Clusters (TAMC VI)", Mexico City, Mexico, June 13-17, 2010

235. Center for High Performance Computing and Communications
Priya Vashishta, "Academic Affairs Committee USC Board of Trustees", University of Southern California, Los Angeles, CA, November 03, 2010

236. Material Properties and Processes Using High Performance Computing
Priya Vashishta, "Supercomputing: The Imperative and the Path Forward", New York University Abu Dhabi Institute, Abu Dhabi, January 9-10, 2011

237. Computational Tools: Current length and time scales Petascales to Exascale Computing
Priya Vashishta, "Workshop on Harnessing DOE's High Performance Computing Expertise to Strengthen the US Chemical Enterprise", Council for Chemical Research and Department of Energy, Washington DC, March 10, 2011

238. Reactive Nanosystems
Priya Vashishta, "3rd International Conference from Nanoparticles and Nanomaterials to Nanodevices and Nanosystems", IC4N-2011, Crete, Greece, June 26, 2011

239. Multimillion to Billion Atom Simulations of Properties and Processes in Materials
Priya Vashishta, "8th International Conference on Diffusion in Materials (DiMat 2011)",

INVITED TALKS AT CONFERENCES:

Dijon, France, July 4-8, 2011

240. Multimillion to Billion Atom Simulations of Properties and Processes in Materials
Priya Vashishta, "High Fidelity 3D Multiscale Materials Modeling and Experimental Analysis Workshop", Vicksburg, MS, August 2–3, 2011

240. Reactive Nanosystems: Multimillion Atom Molecular Dynamics Simulations of Energetic Materials
Priya Vashishta, "Workshop on Nanoenergetics and Combustion Dynamics", Arlington, Virginia, August 23-25, 2011

241. Multimillion to Billion Atom Simulations of Properties and Processes in Materials
Priya Vashishta, "Materials Science and Technology (MS&T 2011)", Columbus, OH, October 14, 2011

242. Multimillion Atom Simulations of Reactive Nanosystems
Priya Vashishta, "The International Workshop on the Physics of Semiconductor Devices", IIT Kanpur, India, December 19-22, 2011

243. Multimillion Atom MolecularDynamics Simulations of Reactive Nanosystems
Priya Vashishta, "PACAM XII", Port of Spain, Trinidad, 2-6 January 2012

244. Shock Compression of Condensed Systems: Role of Emerging Computing Platforms
Priya Vashishta, "Dynamic Compression Sector@APS, Argonne National Laboratory, Argonne, IL, 19-20 January 2012

245. Small Interfering RNA Induces Structural Phase Transformation in a Phospholipid Bilayer
Priya Vashishta, E-MRS Spring Meeting, Strasbourg, France, May 14, 2012

246. Computational Physics and Chemistry of Energetic Materials - Discussion Leader
Priya Vashishta, Gordon Conference on Energetic Materials 2012, Mount Snow resort in Dover, Vermont, June 20, 2012

247. Materials Damage and Self-healing Materials Under Extreme Conditions
Priya Vashishta, DOE-BES Meeting on Theoretical Condensed Matter, Rockville, MD, August 21, 2012

248. Atomistic Simulation of Reactive Materials: Multimillion Atom Molecular Dynamics Simulations of Energetic Materials
Priya Vashishta, Reactive Materials Fundamentals, Synthesis Techniques, and Applications, Materials Research Society, Boston, MA, November 25, 2012

249. Reactive Nanosystems: Multimillion Atom Molecular Dynamics Simulations of Energetic Materials
Priya Vashishta, Workshop on Electronic Structure Approaches to Atoms, Molecules, Clusters and Solids Techniques, and Applications, Hyderabad, India, 7-11 January 2013

250. Multimillion to Billion Atom Simulations of Properties and Processes in Materials
Priya Vashishta, Grand Challenges in Computational Materials Design, North Carolina State University, NC, January 15-16, 2013

251. Reactive Nanosystems: Multimillion Atom Reactive Molecular Dynamics Simulations
Priya Vashishta, CECAM Workshop on Materials chemomechanics at the atomic scale:

INVITED TALKS AT CONFERENCES:

modeling and experiments, Laussane, Switzerland, April 29-may2, 2013

252. Rapid Hydrogen Production from Water Using Aluminum Nanoclusters: A Quantum Molecular Dynamics Simulation Study
Priya Vashishta, Solid State Ionics-19 International Conference, Kyoto, Japan, June 2-7, 2013

253. Atomistic Simulation of Reactive Materials: Multimillion Atom Molecular Dynamics Simulations of Energetic Materials,
Priya Vashishta, Kumamoto University, Kumamoto, Japan, June 6, 2013

254. Self-Healing Materials and Damage From Shock Induced Nanobubble Collapse: Reactive Molecular Dynamics Simulations
Priya Vashishta, Collaborative Conference on 3D & Materials Research (CC3DMR 2014), Incheon/Seoul, South Korea, June 23, 2014

255. Reactive Nanosystems: Multimillion Atom Molecular Dynamics Simulations of Energetic Materials
Priya Vashishta, Daejeon, South Korea, 26 June 2014

256. Thermomechanical Behavior & Materials Damage: Multimillion-Billion Atom Reactive Molecular Dynamics Simulations",
Priya Vashishta, Workshop on High Performance Computing -- From Clouds and Big Data to Exascale and Beyond, Cetraro, Italy, 7-11 July 2014

257. Thermo-mechanical Behavior and Materials Damage : Multimillion-Billion Atom Reactive Molecular Dynamics Simulations,
Priya Vashishta, Mechanical Behavior of Structural Materials, Pan American Materials Conference 2014, Sao Paulo, Brasil, July 24, 2014

258. Self-Healing Nanomaterials and Hydrogen-on-Demand: Multimillion-Atom Reactive Molecular Dynamics Simulations
Priya Vashishta, DOE-Theoretical Condensed Matter Physics Meeting, Materials Sciences and Engineering Division, Rockville, MD, August 11, 2014

259. Multimillion-Billion Atom Reactive Molecular Dynamics Simulations,
Priya Vashishta, EsNet – The DOE Science Network, Germantown, MD, September 10 -11, 2014

260. Self-Healing Materials and Damage From Shock Induced Nanobubble Collapse: Reactive Molecular Dynamics Simulations,
Priya Vashishta, 7th International Conference on Multiscale Materials Modeling, Berkeley, California, October 6-10, 2014

261. Self-Healing Materials and Damage From Shock Induced Nanobubble Collapse: Reactive Molecular Dynamics Simulations
Priya Vashishta, Global Forum on Nanoelectronic Manufacturing: From Materials to Systems, Mumbai, India, October 9, 2014

262. Reactive Molecular Dynamics Simulations, Data Analytics and Visualization
Priya Vashishta, Informatics and Genomics for Materials Development, Materials Research Society, Boston, MA, December 2, 2014

INVITED TALKS AT CONFERENCES:

262. Massively Parallel Reactive and Quantum Molecular Dynamics Simulations
Priya Vashishta, Fifty Years of Molecular Dynamics Simulations I: Past, Present and Future, American Physical Society, Arlington, TX, March 2, 2015

262. Massively Parallel Reactive and Quantum Molecular Dynamics Simulations
Priya Vashishta, RE^3 Workshop 2015: Renewable Energy & Efficiency, Louiville, KY, March 23, 2015

262. Massively Parallel Reactive and Quantum Molecular Dynamics Simulations
Priya Vashishta, 4th International Symposium on Energy Challenges & Mechanics – Working on Small Scales, Aberdeen, Scotland, August 11-13, 2015

263. Self-Healing Materials and Damage From Shock Induced Nanobubble Collapse: Reactive Molecular Dynamics Simulations
Priya Vashishta, Plasticity, Damage and Fracture 2016, Hawaii, January 3-9, 2016

264. Massively Parallel Reactive and Quantum Molecular Dynamics Simulations
Fifty Years of Molecular Dynamics Simulations: Past, Present and Future
Priya Vashishta, American Chemical Society, Computers in Chemistry, San Diego, CA, March 13, 2016

265. Reactive Nanosystems: Billion Atom Reactive and Quantum Molecular Dynamics Simulations
Priya Vashishta, *THERMEC'2016: 9th International Conference on Processing & Manufacturing of Advanced Materials*, Graz, Austria, May 30, 2016

266. Reactive Molecular Dynamics Simulations, Data Analytics and Visualization
Priya Vashishta, International Union of Materials Research Societies (IUMRS2016) International Conference on Electronic Materials (ICEM2016), Symposium on Materials by Theoretical-Computational Design, Singapore, July 4-8, 2016

267. Materials Genome Innovation for Computational Software
Priya Vashishta, Theoretical Condensed Matter Physics, Materials Science & Engineering, Basic Energy Sciences, DOE, Gaithersburg, MD, August 16, 2016

268. Reactive Nanosystems: Billion Atom Reactive and Quantum Molecular Dynamics Simulations
Priya Vashishta, American Chemical Society, Molecular Dynamics for Modern Materials with LAMMPS, Temple University, August 19, 2016

269. Reactive Nanosystems: Billion Atom Reactive and Quantum Molecular Dynamics Simulations
Priya Vashishta, European Materials Research Society (E-MRS), Warsaw, Poland, September 19, 2016

270. Reactive Nanosystems Under Thermomechanical Extremes: Billion Atom Reactive and Quantum Molecular Dynamics Simulations
Priya Vashishta, Materials Research Society (MRS), Boston, November 30, 2016

271. Reactive Nanosystems: Billion Atom Reactive and Quantum Molecular Dynamics Simulations
Priya Vashishta, International Conference on Plasticity, Damage & Fracture, Symposium on Structural Changes in Materials, Puerto Vallarta, Mexico, January 3-9, 2017

INVITED TALKS AT CONFERENCES:

272. Reactive Nanosystems: Billion Atom Reactive and Quantum Molecular Dynamics Simulations Priya Vashishta, International Conference on Computational Physics (ICCP10), Materials for energy storage and conversion on Structural Changes in Materials, Puerto Vallarta, Mexico, January16-20, 2017

INVITED TALKS GIVEN AT UNIVERSITIES AND RESEARCH ORGANIZATIONS:

St. Andrews University, U. K. (1968)

McMaster University, Canada (1970, 1983)

Dalhousie University, Canada (1970)

Northwestern University (1971)

Iowa State University (1972)

Michigan State University (1972, 1994)

IBM Thomas J. Watson Research Center- Yorktown (1972)

Argonne National Laboratory (1972, 1974, 1978, 1981, 1985, 1989. 1991, 1994)

Western Michigan University (1973, 1979)

University of Messina, Italy (1973, 1973)

Roorkee University, India (1973)

Tata Institute of Fundamental Research, India (1973)

Northern Illinois University (1974, 2005)

University of Iowa (1974)

University of Chicago (1974, 1982)

University of Cincinnati (1974)

University of Illinois at Chicago (1974, 1982)

University of California - Berkeley (1975, 1982)

University of British Columbia, Canada (1975, 1983)

Osaka University, Japan (1975, 1985)

Nagoya University, Japan (1975, 1975, 1985)

Purdue University (1975, 1978, 1981, 1984)

Bell Laboratory - Holmdel (1976)

INVITED TALKS GIVEN AT UNIVERSITIES AND RESEARCH ORGANIZATIONS:

University of California - San Diego (1977, 1977, 1982, 2011)

University of California - Los Angeles (1977, 2004)

University of California - Davis (1977, 2004, 2011)

CECAM - Orsay, France (1977)

University of Missouri - Columbia (1978, 1983, 1984)

University of Illinois - Urbana (1979)

Chalmers Institute of Technology, Sweden (1981, 1981)

Western Illinois University (1981)

Arizona State University (1981)

Bell Laboratories - Murray Hill (1981)

Brown University (1981)

Cornell University (1981, 1987)

IBM Research Laboratory - San Jose (1982)

Washington University - St. Louis (1982, 1984, 1991)

California Institute of Technology (1982)

University of Minnesota (1982, 1988)

University of California - Irvine (1982, 2011)

University of Southern California (1982, 2001, 2003)

University of Toronto - Canada (1983)

Simon Fraser University - Canada (1983)

Institute for Theoretical Physics - Santa Barbara (1983, 1994)

University of Houston, Houston, TX (1983, 1996))

Kansas State University (1984)

INVITED TALKS GIVEN AT UNIVERSITIES AND RESEARCH ORGANIZATIONS:

International Center for Theoretical Physics - Trieste, Italy (1984)

Schlumberger-Doll Research Laboratory (1984)

Institute of Solid State Physics, Tokyo, Japan (1985, 1989)

Tokyo University, Japan (1985, 1989)

Kyoto University, Japan (1985, 1989)

Niigata University, Japan (1985, 1989, 2007)

Tohoku University, Japan (1985, 1989)

University of Sao Paulo, Sao Carlos, Brazil (1985, 1987)

Louisiana State University (1988, 1989)

Australian National University, Canberra, Australia (1989)

Okayama University, Japan (1989, 1995)

Kanazawa University, Japan (1989)

Clemson University (1990)

Punjab University, Chandigarh, India (1991)

National Physical Laboratory, Delhi, India (1991)

Institute of Physics - Bhubaneswar, India (1991)

Indian Institute of Science - Bangalore, India (1991)

Emory University (1991)

Texas A&M University (1991)

Minnesota Supercomputing Research Institute (1991)

NEC Corporation, Tsukuba, Japan (1992)

Government Industrial Research Institute, M I T I, Osaka, Japan (1992)

University of New South Wales, Sydney, Australia (1993)

INVITED TALKS GIVEN AT UNIVERSITIES AND RESEARCH ORGANIZATIONS:

Institute of Advanced Study, Canberra University, Canberra, Australia (1993)

National Institute for Standards and Technology (NIST), Washington DC (1994)

Energy Research Laboratory, Hitachi City, Japan (1994)

Argonne National Laboratory, Argonne, IL (1996)

Northwestern University, Evanston, IL (1996)

Institute of Theoretical Physics, University of California, Santa Barbara (1997)

Argonne National Laboratory, Argonne, IL (1997)

NASA Ames Laboratory, CA (1997)

Ohio State University, Columbus, OH (1997)

Rice University, Houston, TX (1998)

Ohio State University (1998)

Cornell University, Ithaca, NY (1999)

Florida State University, Tallahassee, FL (2000)

Harvard University, Cambridge, MA (2000, 2008)

Universidade Federal de São Carlos, São Carlos, Brasil (2000)

Universidade Federal (UFAL), Maceio, Brasil (2000)

North Carolina State University, NC (2000)

Livermore National Laboratory, CA (2001)

General Motors, MI (2001)

Fermi National Laboratory, IL (2001)

University of Georgia, GA (2001)

Sandia National Laboratory (2001)

Agilent Technologies, CA (2003)

Army Research Laboratory, Aberdeen Proving Grounds, MD (2004)

University of Pittsburgh, PA (2007)

University of Uppsala, Uppsala, Sweden (2009)

University of South Florida, Tampa, Florida (2011)

INVITED TALKS GIVEN AT UNIVERSITIES AND RESEARCH ORGANIZATIONS:

University of Central Florida, Orlando, Florida, September (2013)

Louisiana State University, Baton Rouge, Louisiana (2014)

University of Connecticut, Storr, Connecticut (2015)