



Curriculum Vitae	
Personal information	
First name(s) / Surname(s)	Hiqmet Kamberaj
Address(es)	International Balkan University, Tashko Karadza 11A, 1000, Skopje, R. of Macedonia
Telephone(s)	+389(0)23214831 (ext. 123)
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E-mail	hkamberaj@ibu.edu.mk or h.kamberaj@gmail.com
Web address	hkamberajibu.wikidot.com
Nationality	Albanian, Macedonian
Date of birth	07/09/1972
Education	<ol style="list-style-type: none"> 1. Ph.D. in Physics, 2001-2005, Manchester Metropolitan University, Manchester, United Kingdom. 2. M.Sc. in Physics, 1997-2000, University of Siegen, Siegen, Germany. 3. B.Sc. (5 years study program), 1991-1996, University of Tirana, Tirana, Albania.
Languages Native Foreign languages (Reading/ writing/ speaking)	<ol style="list-style-type: none"> 1. Albanian 2. English, German
h-Index Google Scholar ResearchGate Mendeley	9 to 10 https://scholar.google.com/citations?user=5I1XAAcAAAAJ&hl=en https://www.researchgate.net/profile/Hiqmet_Kamberaj https://www.mendeley.com/profiles/hiqmet-kamberaj2/?utm_source=mm_Mandrill

<p>ACADEMIC EXPERIENCE</p> <p>Visiting scholar</p> <p>Postdoctoral positions</p> <p>Graduate student positions</p> <p>Teaching positions</p>	<ol style="list-style-type: none"> 1. 2014, Visiting Scholar Researcher, University of South Florida, USA 2. 2008-2009, Postdoctoral Research Associate, Center for Drug Design, University of Minnesota, MN, USA. 3. 2006-2008, Postdoctoral Research Associate. Center for Biological Physics, Arizona State University, Department of Chemistry & Biochemistry, Tempe, AZ, USA. 4. 2005-2006, Postdoctoral Fellow, University of Alberta, National Institute of Nanotechnology, Edmonton, Alberta, Canada. 5. 2004-2005, Postdoctoral Research Associate, Manchester Metropolitan University, Manchester, United Kingdom. 6. 2000-2001, Postgraduate Student, Max-Planck Institute of Biophysics, Frankfurt am Main, Germany. 7. 2014-Present, Associate Professor, International Balkan University, Skopje, Republic of Macedonia. 8. 2014-Present, Adjunct Associate Professor, State University of Tetovo, Tetovo, Republic of Macedonia. 9. 2009-2014, Assistant Professor, International Balkan University, Skopje, Republic of Macedonia. 10. 2011-2014, Adjunct Assistant Professor, State University of Tetovo, Tetovo, Republic of Macedonia. 11. 2004-2005, Assistant Lecturer, Manchester Metropolitan University, Manchester, United Kingdom. 12. 2003-2004, Teaching Assistant, Manchester Metropolitan University, Manchester, United Kingdom. 13. 2001-2003, Teaching Assistant, University of Coventry, Coventry, United Kingdom. 14. 1996-1997, Assistant Lecturer, University of Tirana, Tirana, Albania.
<p>Teaching Assignments</p> <p>Courses at IBU</p>	<p><u>Undergraduate Level:</u></p> <ul style="list-style-type: none"> • Introduction to Programming • Computer Architecture • Discrete Computational Structure • Discrete Mathematics • Algorithms • Systems Modeling • Introduction to Information Technology • Introduction to Probability & Statistics • Physics I • Physics II • Electromagnetic Theory

<p>Other Universities (State University of Tetovo)</p>	<p><u>Graduate Level:</u></p> <ul style="list-style-type: none"> • Information Theory • Theory of Probability <p><u>Undergraduate Level:</u></p> <ul style="list-style-type: none"> • Statistical Physics • Classical Mechanics • Electrodynamics • Medical Physics <p><u>Graduate Level:</u></p> <ul style="list-style-type: none"> • Theoretical Methods in Physics
<p>Research interests and fields</p>	<ul style="list-style-type: none"> • Theoretical and Computational Physics/Chemistry • Soft matter, Biophysics and Nanophysics • Protein and DNA dynamics • Bio-inspired optimization • Bioinformatics • Numerical integrators for molecular dynamics • Monte Carlo and molecular dynamics techniques • Free energy calculations • Information flow in biomolecular systems • Coarse-grained models of proteins and DNA • Swarm particle intelligence optimization • GPU-accelerated computing • Quantum simulations of molecular systems
<p>Administrative Working Experience</p> <p>Professional Memberships</p>	<ul style="list-style-type: none"> • Acting Dean of Faculty of Engineering (2017-Present): International Balkan University • Rector's Advisor for Research and Projects (2015-Present): International Balkan University • President of Student Evaluation Committee (2012-Present): International Balkan University • Member of the Senate (2009-Present): International Balkan University • Member of the Faculty Council (2009-Present): International Balkan University <ul style="list-style-type: none"> • 2014-Present Research Institute of Science and Technology (Albania) • 2017-Present Subscriber Member of Nature Methods Techniques for Life Scientists and Chemists (USA) • 2011-2016 American Chemical Society (USA) • 2006-2009 American Biophysics Society (USA) • 2001-2005 British Liquid Crystal Society (UK)

Professional Experience	<ul style="list-style-type: none"> • Editor-in-Chief of American Journal of Applied Sciences, Engineering and Technology (2014-2015). • Editorial Board Member of: International Journal of Dataset Papers in Physics (2012-2017); Turkish Academic Journal of Intelligent Systems. • Peer Reviewer of: International Journal of Computational and Theoretical Chemistry; <u>International Journal of Electrical Power & Energy Systems</u>; Science Journal of Applied Mathematics and Statistics; Journal of Chemical Information and Modeling; BMC Bioinformatics; Entropy.
Awards	<ol style="list-style-type: none"> 1. 2016: Certificate of Excellence (Motivation: For his outstanding performance in academic research during the academic year 2015/2016), International Balkan University. 2. First Prize in Bioinformatics: Building Bridges Symposium, Minneapolis, MN, USA (17 April 2009). Work presented as a poster – <i>Computational Alanine scanning of β-lactamase Inhibitor Protein and TEM-1 β-lactamase Complex</i>. 3. Summer 1996: Dean’s list of outstanding students, University of Tirana, Tirana, Albania. 4. April 1986: Second prize of the mathematical Olympiad, Vlora, Albania.
PUBLICATIONS	
Original scientific article (Including electronic journals) (2014-Present)	<ol style="list-style-type: none"> 1. H. Kamberaj, (2018) Faster Protein Folding Using Enhanced Conformational Sampling of Molecular Dynamics Simulation, <i>Journal of Molecular Graphics and Modelling</i>, 81, 32-49. 2. R. Izairi and H. Kamberaj, (2017) Comparison Study of Polar and Non-polar Contributions to Solvation Free Energy, <i>Journal of Chemical Information and Modeling</i>, 57(10), 2539-2553. 3. H. Kamberaj, (2017) Sampling Convergence of Collective Motions in Proteins, <i>Journal of Applied Physical Science International</i>, 8(3), 101-112. 4. H. Kamberaj, (2016) Replica Exchange Using q-Gaussian Swarm Quantum Particle Intelligence Method, <i>Engineering and Applied Sciences</i>, 1(2), 20-25. 5. H. Kamberaj, (2015) Conformational Sampling Enhancement of Replica Exchange Molecular Dynamics Simulations Using Swarm Particle Intelligence, <i>The Journal of Chemical Physics</i>, 143, 124105. 6. H. Kamberaj, (2014) Q-Gaussian Swarm Quantum Particle Intelligence on Predicting Global Minimum of Potential Energy Function, <i>Applied Mathematics and Computation</i>, 229, 94-106. <ol style="list-style-type: none"> 7. Spiriti, J., Kamberaj, H., de Graff, A., Thorpe, M.P., and van der Vaart, A., (2012) DNA bending through large angles is aided by ionic screening, <i>Journal of Chemical Theory and Computation</i>, 8 (6), 2145–2156. 8. Spiriti, J., Kamberaj, H., and van der Vaart, A., (2012) Development and application of enhanced sampling techniques to simulate the long-time scale dynamics of biomolecular systems, <i>International Journal of Quantum Chemistry</i>, 112, 33-43. 9. Kamberaj, H. (2011) A Theoretical Model for the Collective Motion of Proteins by Means
(2009-2013)	

<p style="text-align: center;">(2001-2008)</p>	<p>of Principal Component Analysis. <i>Central European Journal of Physics</i>, 9(1), 96-109.</p> <p>10. Kamberaj, H. and van der Vaart, A. (2009) Extracting the causality of correlated motions from molecular dynamics simulations. <i>Biophysical Journal</i>, 97(6), 1747-1755.</p> <p>11. Kamberaj, H. and van der Vaart, A. (2009) An optimized replica exchange method for molecular dynamics simulations. <i>Journal of Chemical Physics</i>, 130(7), 074904-6.</p> <p>12. Kamberaj, H. and van der Vaart, A. (2009) Correlated Motions and Interactions at the Onset of the DNA-Induced Partial Unfolding of Ets-1. <i>Biophysical Journal</i>, 96, 1307-1317.</p> <p>13. Kamberaj, H. and van der Vaart, A. (2007) Multiple scaling replica exchange method for the efficient sampling of biomolecular systems. <i>The Journal of Chemical Physics</i>, 127(23), 234102-7.</p> <p>14. Kamberaj, H., Low, R.J., and Neal, M.P. (2006) Correlation between molecular chirality and helical twisting power. A computer simulation study. <i>Molecular Physics</i>, 104(3), 335-357.</p> <p>15. Neal, M.P., Kamberaj, H., and Low, R.J. (2005) Some calculations of molecular chirality. <i>Molecular Crystal Liquid Crystal</i>, 439, 1937-1943.</p> <p>16. Kamberaj, H., Low, R.J., and Neal, M.P. (2005) Molecular structure and helical twisting power, <i>Ferroelectrics</i>, 315, 183-196.</p> <p>17. Kamberaj, H., Low, R.J., and Neal, M.P. (2005) Symplectic and time reversible integrators for molecular dynamics simulations of rigid molecules. <i>The Journal Chemical Physics</i>, 122(22), 224114-30.</p> <p>18. Kamberaj, H., Osipov, M.A., Low, R.J., and Neal, M.P. (2004) Helical twisting power and chirality indices. <i>Molecular Physics</i>, 102(5), 431-446.</p> <p>19. Kamberaj, H., Low, R.J., and Neal, M.P. (2003) Symplectic and time reversible integrators for rigid bodies. <i>System Science</i>, 29(2), 31-45.</p> <p>20. Kamberaj, H., and Helms, V. (2001) Monte Carlo simulation of biomolecular systems with BIOMCSIM, <i>Computer Physics Communications</i>, 141 (3), 375-402.</p>
<p style="text-align: center;">Published scientific conference contribution Full paper</p>	<p>1. Kamberaj, H., Low, R.J., and Neal, M.P. (2003) <i>Symplectic and time reversible integrators for rigid bodies</i>, 16th Conference on Systems Engineering, ICSE2003, editors: K.J. Burnham and O.C.L. Haas, Coventry University, Coventry, UK, Vol. 1, 320-325, September 9-11.</p> <p>2. T. Conka-Nurdan, K. Nurdan, F. Constantinescu, B. Freisleben, H. Kamberaj, N.A. Pavel, I. Rauhut, K. Reichmann, A.H. Walenta (2001) <i>Influence of the Detector Parameters on a Compton Camera</i>, IEEE, 22-26.</p>

<p>Published scientific conference contribution abstract Abstract</p>	
<p>(2014-Present)</p>	<ol style="list-style-type: none"> 1. Izairi-Bexheti, R. and Kamberaj, H., On the calculation protein-protein binding free energy using molecular dynamics simulations, 1st International Conference of Natural Sciences and Mathematics, Tetovo, R. of Macedonia, June 16-17, 2017. 2. Kamberaj, H., Perspectives on Molecular Dynamics Simulations & Free Energy Calculations for Biomolecular Systems, International Conference on Applied Sciences, Engineering and Mathematics, Ohrid, R. of Macedonia, May 5-7, 2017. 3. Izairi, R. and Kamberaj, H., <i>Molecular Dynamics Simulation Study of Protein-Protein Interactions</i>, ALBNANO2016 NANOTECHNOLOGY AND BIOSENSORS, May 29-31, 2016, Tirana, Albania. 4. Kamberaj, H., 2nd International Conference on Pure & Applied Science, June 1-5 2016, Istanbul Turkey, <i>Enhanced Transition Path Sampling Using Swarm Particle Intelligence in Replica Exchange Molecular Dynamics</i>. 5. Izairi, R. and Kamberaj, H., The 10th Conference of the Society of Physicists of Macedonia, Skopje, Macedonia, September 25-28, 2014, <i>Calculations of the Electronic Properties of Molecules using Computational Quantum Mechanics Models</i>. 6. Mutlu, O., Kamberaj, H., 2nd International Bau-Drug Design Symposium, <i>Novel methods and emerging targets in drug discovery & patented drug development</i>, Bahcesehir University, Istanbul, Turkey, 17-19 April 2014, <i>Structural Modelling, Molecular Docking and Molecular Dynamics Simulation of Trypanothione Synthase</i>.
<p>(2009-2013)</p>	<ol style="list-style-type: none"> 1. Kamberaj, H., International Conference on Applied Analysis and Mathematical Modeling, Istanbul, Turkey, 2-5 June 2013, <i>Symbolic Mutual Information for Estimation of Nonlinear Higher Order Correlated Fluctuations</i>. 2. Kamberaj, H., International Meeting on Atomic and Molecular Physics and Chemistry, Scuola Normale Superiore Pisa, Italy (12-14 September 2012), <i>Dynamics and static dielectric response of water in anionic hydration shells using ab-initio molecular dynamics simulation</i>. 3. Kamberaj, H., 6th International Annual Meeting of Alb-Science Institute, Prishtina, Kosova (1-4 September 2011), <i>Modeling of protein-protein interactions: Calculation of the binding free energy and computational alanine scanning</i>. 4. Kamberaj, H., 8th European Biophysics Congress, Budapest, Hungary (23-27 August 2011), <i>Elucidating the modular structure of the protein G C2 fragment and human IgG Fc domain binding site using computer simulations</i>.

(2002-2008)	<ol style="list-style-type: none"> 1. Kamberaj, H. and van der Vaart, A. (2008) Biophysics Society 52-th Annual Meeting & 16-th IUPAB International Biophysics Congress, Long Beach, CA, USA. Work presented as a poster - <i>Unfolding upon binding: Elucidation of the complex binding dynamics of the Ets-1 transcription factor by computer simulation</i>, February 2-6. 2. Neal, M.P., Kamberaj, H. and Low, R.J. (2006) 21st International Liquid Crystal Liquid Crystals Conference 2006, Keystone, Colorado, USA. Work presented as poster – <i>On predicting ferroelectric liquid crystal reduced spontaneous polarization or polarization power</i>, July 2-7. 3. Kamberaj, H., Low, R.J., and Neal, M.P. (2005) 6th International Conference on Liquid Matter at Utrecht University, Utrecht, The Netherlands. Work presented as a poster – <i>Calculation of molecular chirality and helical twisting power</i>, July 1-6. 4. Kamberaj, H., Low, R.J., and Neal, M.P. (2005) British Liquid Crystal Society Annual Meeting at Exeter University, Exeter, UK. Work presented as a talk - <i>Calculations of the Molecular Chirality</i>, March 22-26. 5. Kamberaj, H., Low, R.J., and Neal, M.P. (2004) British Liquid Crystal Society Annual Meeting at Manchester Metropolitan University, Manchester, UK. Work presented as a talk and in poster format - <i>Applications of Calculations of the Molecular Chirality and Helical Twisting Power</i>, April 5-7. 6. Kamberaj, H., Low, R.J. and Neal, M.P. (2003) 9th International Conference on Ferroelectric Liquid Crystals at Trinity College, Dublin, Ireland. Work presented in poster format – <i>Molecular structure and helical twisting power</i>, August 24-29. 7. Kamberaj, H., Low, R.J., and Neal, M.P. (2003) British Liquid Crystal Society Annual Meeting at Cambridge University, Cambridge, UK. Work presented in poster format, April 7-9. 8. Kamberaj, H., Low, R.J., and Neal, M.P. (2002) 19th International Liquid Crystal Conference 2002, Edinburgh, UK. Work presented in poster format, June 30-July 5.
Published professional conference contribution abstract	<ol style="list-style-type: none"> 1. Kamberaj, H. (2011) Elucidating the modular structure of the protein G C2 fragment and human IgG Fc domain binding site using computer simulations, <i>European Biophysics Journal with Biophysics Letters</i>, 40, 108-108. 2. van der Vaart, A., Kamberaj, H. (2009) Extracting The Causality Of Correlated Motions From Molecular Dynamics Simulations, <i>Biophysical Journal Abstracts</i>, 96(3), pp 406a.
MONOGRAPHS AND OTHER COMPLETED WORKS	

<p>Textbook (M.Sc. and Ph.D. graduates, and researchers level):</p> <p>Textbook (undergraduate level):</p>	<ol style="list-style-type: none"> 1. Kamberaj, H. (2016) <i>Molecular Dynamics Simulations in Statistical Physics. Theory and Applications</i>. State University of Tetovo. 1. Kamberaj, H. (2014) <i>Probability and Statistics. Essentials about Probability and Statistics</i>, International Balkan University. Skopje, R. of Macedonia. 2. Kamberaj, H. (2014) <i>Thermodynamics & Statistical Physics</i>, State University of Tetovo, Tetovo, R. of Macedonia. 3. Kamberaj, H. (2012) <i>Fundamentals of Medical Physics</i>, State University of Tetovo, Tetovo, R. of Macedonia. 4. Kamberaj, H. (2010) <i>Introduction to Programming. An undergraduate engineering student textbook</i>, International Balkan University, Skopje, R. of Macedonia.
<p>Doctoral dissertation</p>	<ol style="list-style-type: none"> 1. Kamberaj, H. (2005) <i>Simulation and modelling of liquid crystal mesophases linked to the design of real materials</i>, Manchester Metropolitan University, Manchester, United Kingdom, Ph.D. Thesis. Supervisors: Prof. Dr. Maureen Neal and Prof. Dr. Robert Low
<p>Master's thesis</p>	<ol style="list-style-type: none"> 1. Kamberaj, H. (2000) <i>The Influence of Detector Parameters on the Image Reconstruction for a Compton Camera</i>, University of Siegen, Siegen, Germany. Master Thesis. Supervisor: Prof. Dr. Nikolae Pavel
<p>Undergraduate thesis</p>	<ol style="list-style-type: none"> 1. Kamberaj, H. (1996) <i>Thermodynamics and Dynamics of Two Dimensional Ising Spin Ladder Lattice Model</i>, B.Sc. Thesis (in Albanian), University of Tirana, Tirana, Albania. Supervisor: Prof. Dr. Rexhep Meidani
<p>Final research report</p>	<ol style="list-style-type: none"> 1. Kamberaj, H. (2013) <i>Symbolic Mutual Information for Estimation of Nonlinear High Order Correlated Fluctuations</i>, International Balkan University, Skopje, Macedonia. 2. Kamberaj, H. (2012) <i>Sampling Convergence of Collective Motions of Proteins</i>, International Balkan University, Skopje, Macedonia. 3. Kamberaj, H. (2006) <i>Dynamics of the Proteins and Principal Component Analysis</i>, University of Alberta, Edmonton, Canada. 4. Kamberaj, H. and Helms, H. (2001) <i>Thermodynamics of internal water molecules in the protein BPTI studied by Monte Carlo simulations</i>, Max-Planck Institute of Biophysics, Frankfurt am Main, Germany.
<p>Software</p>	<ol style="list-style-type: none"> 1. MMGBSA (in Fortran 90), A software for calculation of the solvation free energy of (bio)molecular systems (including polar and nonpolar terms). 2. BIOMCSIM (in C), A Monte Carlo software for computer simulation of biomolecular systems (Published in Computer Physics Communication journal). 3. SQPO (in C++), A Swarm Quantum Optimization general purpose code. 4. MDGOSim (in Fortran, C++), Swarm Particle Replica Exchange Molecular dynamics simulations of proteins using coarse-grained methods. 5. MDAnalTools (Matlab, Fortran, Python), A collection of programs and scripts for modeling and statistical analyzing of molecular dynamics simulations.
<p>PERFORMED WORKS (EVENTS)</p>	
<p>Invited lecture at foreign university</p>	<ol style="list-style-type: none"> 1. Kamberaj, H. (2010) Causality of Correlated Motions in Biomolecular Systems & Novel Algorithm for Global Optimization Problems in Ligand and/or peptide-protein binding, Schroedinger Company, New York, NY USA, November 16.

<p>Unpublished conference contribution</p>	<ol style="list-style-type: none"> 1. Kamberaj, H., and Sham, Y. (2009) Bioinformatics: Building Bridges Symposium, Minneapolis, MN, USA. Work presented as a poster – Computational Alanine Scanning of β-lactamase Inhibitor Protein and TEM-1 β-lactamase Complex, April 17. 2. Kamberaj, H. and Sham, Y. (2009) MIKI Meeting, Minneapolis, MN, USA. Work presented as a poster – Computational Alanine Scanning of β-lactamase Inhibitor Protein and TEM-1 β-lactamase Complex, April 3-5. 3. Kamberaj, H. and van der Vaart, A. (2007) Postdoctoral Research Day symposium at Biodesign Center, Tempe, Arizona, USA. Work presented as a poster - Unfolding upon binding: Elucidation of the complex binding dynamics of the Ets-1 transcription factor by computer simulation, October 6. 4. Kamberaj, H. (2007) Annual meeting BioPhest 2007 at Arizona State University, Tempe, Arizona, USA. Work presented as a talk - Multiple scaling replica exchange method for the efficient sampling of biomolecular systems, April 21. 5. Kamberaj, H. (2006) Postdoctoral Research Day Symposium at University of Alberta, Edmonton, Alberta, Canada. Work presented as a poster – Multiple time scale molecular dynamics simulations of biomolecular systems, April 7.
<p>Unpublished invited conference lecture</p>	<ol style="list-style-type: none"> 1. Kamberaj, H. (2010) 8th Conference of the Society of Physicists of Republic of Macedonia, Skopje, R. of Macedonia. The role of the correlated motions in biomolecular systems, September 23-25. 2. Kamberaj, H. and Sham, Y. (2009) Gordon Conference on Computer aided Drug Design, Tilton, NH, USA. Computational Alanine Scanning of β-lactamase Inhibitor Protein and TEM-1 β-lactamase Complex, July 19-24. 3. Kamberaj, H. (2004) Mini-Symposium on the Chirality and Nonlinearity in Liquid Crystals, The Rank Prize, UK. Work presented as a talk - Calculations of Molecular Chirality, November 29-December 2. 4. Kamberaj, H. (2004) Young Materials Modellers Forum at Daresbury Laboratory, UK. Work presented as a talk - Symplectic and Time Reversible Integrators for Molecular Dynamics Simulations of Rigid Molecules, May 27.
<p>SCIENTIFIC RESEARCH PROJECTS</p>	

Principal Investigator:	<ol style="list-style-type: none"> 1. Kamberaj, H. (2014-Present) Thermodynamics of biomolecular systems – Free energy calculations, International Balkan University, Skopje, R. of Macedonia. 2. Kamberaj, H. (2012-Present) <i>Bio-inspired optimization methods</i>, International Balkan University, Skopje, R. of Macedonia. 3. Kamberaj, H. (2012-Present) <i>Dielectric response functions of solvent and proteins</i>, International Balkan University, Skopje, R. of Macedonia. 4. Kamberaj, H. (2011-Present) <i>Dynamics and Static Dielectric Response of Water in ionic hydration shells using quantum mechanics simulations</i>, International Balkan University, Skopje, R. of Macedonia. 5. Kamberaj, H. (2010-Present) <i>Protein-ligand and Protein-peptide flexible docking using Swarm Intelligence</i>, International Balkan University, Skopje, R. of Macedonia. 6. Kamberaj, H. (2010-Present) <i>Characterization of the dynamics and thermodynamics of the protein-protein interface. A case study of IgG protein and C2 fragment of protein G</i>, International Balkan University, Skopje, R. of Macedonia. 7. Kamberaj, H. (2009-2010) <i>Collective motions of proteins by means of principal component analysis</i>, International Balkan University, Skopje, R. of Macedonia.
Team Member: (2009-Present)	<ol style="list-style-type: none"> 1. Kamberaj, H. and Mutlu, O., (2016-Present), <i>The protein-protein interaction of HDAC1/2 and CK2 complex</i>, Marmara University Faculty of Science and Arts, Department of Biology 34722, Istanbul, Turkey. 2. Kamberaj, H. and van der Vaart, A., (2014) <i>Implementation of the Transfer Entropy Measures in CHARMM Software</i>, University of South Florida, Tampa, FL, USA. 3. Isler, Y. and Kamberaj, H., (2014-Present) <i>Characterisation of Molecular Systems for Refrigerators</i>, International Balkan University, Skopje, R. of Macedonia. 4. Mutlu, O., Kamberaj, H., (2014) <i>Structural Modelling, Molecular Docking and Molecular Dynamics Simulation of Trypanothione Synthase</i>, International Balkan University, Skopje, R. of Macedonia. 5. Spiriti, J., Kamberaj, H., de Graff, A., Thorpe, M.P., and van der Vaart, A., (2011-2012) <i>DNA bending through large angles is aided by ionic screening</i>, International Balkan University, Skopje, R. of Macedonia. 6. Spiriti, J., Kamberaj, H., de Graff, A., and van der Vaart, A., (2011-2012) <i>Development and application of enhanced sampling techniques to simulate the long-time scale dynamics of biomolecular systems</i>, International Balkan University, Skopje, R. of Macedonia. 7. Kamberaj, H., Eggum, J., Muthyala, R., and Sham, Y. (2008-2009) <i>Vancomycin Resistance: Computational Modeling of Substrate Binding to VanX</i>, Center for Drug Design, University of Minnesota, USA. 8. Kamberaj, H. and Sham, Y. (2008-2009) <i>Designing and docking new inhibitors for HIV protein</i>, Center for Drug Design, University of Minnesota, USA. 9. Kamberaj, H and Sham, Y. (2008-2009) <i>Designing and Docking inhibitors for bacteria TEM-1 protein</i>, Center for Drug Design, University of Minnesota, USA. 10. Kamberaj, H. and Sham, Y. (2008-2009) <i>Modulation of the BPTI-TEM-1 protein-protein interface using computational alanine scanning</i>, Center for Drug Design, University of Minnesota, USA.

<p>(2000-2008)</p>	<ol style="list-style-type: none"> 1. Kamberaj, H. and van der Vaart, A. (2006-2008) <i>Investigating replica exchange method for efficient conformational sampling of biomolecular systems in explicit water</i>, Center for Biological physics, Arizona State University, Tempe, USA. 2. Kamberaj, H. and van der Vaart, A. (2006-2008) <i>Investigating correlated motions for Ets-1 –DNA complex in explicit water</i>, Center for Biological physics, Arizona State University, Tempe, USA. 3. Kamberaj, H. and van der Vaart, A. (2006-2008) <i>Applying non-Boltzmann replica exchange method for studying helix unfolding of Ets-1 upon binding to DNA</i>, Center for Biological physics, Arizona State University, Tempe, USA. 4. Kamberaj, H. and van der Vaart, A. (2006-2008) <i>Extracting causality of correlated motion of the Et-1 protein in apo and bound to DNA states using transfer entropy</i>, Center for Biological physics, Arizona State University, Tempe, USA. 5. Kamberaj, H. and van der Vaart, A. (2006-2008) <i>Implementation of the Hamiltonian replica exchange method in CHARMM program</i>, Center for Biological physics, Arizona State University, Tempe, USA. 6. Kamberaj, H. and Stepanova, M. (2005-2006) <i>Developing coarse-grained models for proteins using principal component analysis</i>, National Institute of nanotechnology, University of Alberta, Edmonton, Canada. 7. Kamberaj, H. and Stepanova, M. (2005-2006) <i>Molecular dynamics simulation of Fc fragment of IgG protein in complex with C2 fragment of protein G using Amber program</i>, National Institute of nanotechnology, University of Alberta, Edmonton, Canada. 8. Kamberaj, H., Low, R.J., and Neal, M.P. (2005) <i>Analyzing the correlation between the scaled chiral indices and the molecular polarization</i>, Manchester Metropolitan University, Manchester, United Kingdom. 9. Kamberaj, H., Low, R.J., and Neal, M.P. (2005) <i>Analyzing the correlation between the scaled chiral indices and the helical twisting power of chiral molecules dissolved in nematic phase</i>, Manchester Metropolitan University, Manchester, United Kingdom. 10. Kamberaj, H. and Helms, V. (2000-2001) <i>Developing grand canonical Monte Carlo method for investigating the concentration of water molecules inside the protein cavities</i>, Max-Planck Institute of Biophysics, Frankfurt am Main, Frankfurt, Germany. 11. Kamberaj, H. and Helms, V. (2000-2001) <i>Characterizing the dynamics and thermodynamics of water molecules inside the protein cavities</i>, Max-Planck Institute of Biophysics, Frankfurt am Main, Frankfurt, Germany. 12. Kamberaj, H. and Pavel, N. (1999-2000) <i>Image reconstruction using the Compton camera. Optimization of detector parameters on image quality</i>, University of Siegen, Siegen, Germany.
<p>MENTORSHIPS</p>	<p>Master Studies:</p> <ol style="list-style-type: none"> 1. Redona Izairi, M.Sc. (2016-2018): Study of Protein-Protein Interactions Using the Molecular Dynamics Simulations, State University of Tetova, Tetova, R. of Macedonia. 2. Granit Nebiu, B.Sc. (2017-2018): The Semantics of Web Design and Development, International Balkan University, Skopje, R. of Macedonia. 3. Dhurata Nebiu, B.Sc. (2017-Present): Local and Average Transfer Entropy as an Information Flow Measure, International Balkan University, Skopje, R. of Macedonia. <p>Bachelor Studies:</p> <ol style="list-style-type: none"> 1. Donika Rexhepi, Fundamental Concepts of Classical Cryptosystems, Department of In-

	<p>formation Technology, International Balkan University, Skopje, R. of Macedonia, 2017.</p> <ol style="list-style-type: none"> 2. Haris Zenovic, Design and development of a web-based platform for teaching and learning, Department of Information Technology, International Balkan University, Skopje, R. of Macedonia, 2017. 3. Festim Kamberi Optimized University Students and Courses Information System, Department of Information Technology, International Balkan University, Skopje, R. of Macedonia, 2017. 4. Argjent Halili, Graphical User Interface for Data Analysis with Visual Molecular Dynamics Program using Tcl/Tk, Department of Information Technology, International Balkan University, Skopje, R. of Macedonia, 2016. 5. Skofiar Kamberi, Design and Application of a Database for Students, Department of Information Technology, International Balkan University, Skopje, R. of Macedonia, 2015. 6. Redona Izairi, Calculations of the Electronic Properties of Molecules using Computational Quantum Mechanics Models, State University of Tetovo, Tetovo 2014.
<p>INVITED LECTURES TRAINING AND STAYS Training</p>	<ol style="list-style-type: none"> 1. August –September 2014, Visiting Researcher in Computational Chemistry, University of South Florida, Department of Chemistry, Tampa, FL, USA. 2. January 2002, Workshop at Manchester Research centre for computational science on Visualization and virtual reality for scientific applications, Manchester, United Kingdom. 3. October 2001, Autumn School at Manchester University on High performance computing and SGI NUMAflex, Manchester, United Kingdom. 4. December 2001, British Liquid Crystal Society Winter Workshop at Hull University, Hull, United Kingdom. 5. June 2000, Summer school at UMIST Manchester on Molecular modeling and computer simulations, Manchester, United Kingdom.
<p>FELLOWSHIP/ SCHOLARSHIP</p>	<ol style="list-style-type: none"> 1. Oct 1997-Feb 2000, Studentship awarded of graduate studies by German Academic of Science, University of Siegen, Siegen, Germany. 2. Feb 2000–Aug 2001, Studentship awarded of postgraduate studies by Deutsche Forschungsgemeinschaft at Max-Planck Institute of Biophysics, Frankfurt/Main, Germany. 3. Sept 2001–Aug 2003, Scholarship awarded of postgraduate studies by Coventry University, Coventry, United Kingdom. Pursued the M.Phil. in Physics. 4. Sept 2003-Dec 2004, Scholarship awarded of postgraduate studies by Manchester Metropolitan University, Manchester, United Kingdom. Pursued the Ph.D. in Physics. 5. Jan 2006-Dec 2006, Postdoctoral Fellowship awarded of postdoctoral studies by Alberta University, Edmonton, Canada.