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Obrazovanje

- 1992. diplomirala na Fakultetu za fizičku hemiju, Univerzitet u Beogradu; tema: Primena metode DVR na računanje vibracionih stanja nekih molekularnih sistema
- 1995. magistar fizičko-hemijskih nauka; tema: *Ab initio* računanje spektra radikala BH₂, Fakultet za fizičku hemiju, Univerzitet u Beogradu
- 1997. doktor fizičko-hemijskih nauka; tema: *Ab initio* proučavanje vibranske sprege u četvoratomskim molekulima – analiza spektra B₂H₂, Fakultet za fizičku hemiju, Univerzitet u Beogradu

Profesionalno iskustvo

- 1992 – 2000: asistent na Fakultetu za fizičku hemiju, Univerzitet u Beogradu
- 2006 – 2007: istraživač-saradnik, Institut za hemiju, tehnologiju i metalurgiju, Univerzitet u Beogradu
- 2007 - 2012: naučni saradnik, Institut za hemiju, tehnologiju i metalurgiju, Univerzitet u Beogradu
- 2012 - : viši naučni saradnik, Institut za hemiju, tehnologiju i metalurgiju, Univerzitet u Beogradu

Boravci u inostranstvu

- 1995 Institut za fizičku i teorijsku hemiju, Univerzitet u Bonu, Bon, Nemačka
- 1999 – 2000: Institut za teorijsku hemiju, Univerzitet u Dizeldorfu, Dizeldorf, Nemačka
- 2000 – 2001: Steacie institut za molekularne nauke, Otava, Kanada

Nagrade i priznanja

- Nagrada za najboljeg studenta generacije koji je na Fakultetu za fizičku hemiju diplomirao 1992. godine, Univerzitet u Beogradu
- Nagrada za najbolji diplomski rad odbranjen 1992. godine na Fakultetu za fizičku hemiju, Univerzitet u Beogradu

Stipendije

- Stipendija za postdoktorske studije, Alexander von Humboldt fondacija, Nemačka
- Stipendija za postdoktorske studije, NSERC, Kanada

Oblasti istraživanja

- strukturne, vibracione i elektronske osobine polutanata i molekula od biohemijskog značaja
- računanje strukture i spektara manjih poliatomskih molekula
- ispitivanje uloge supstituenata (alkil, nitro, halogeni itd.) policikličnih aromatičnih ugljovodonika na molekulske osobine koje mogu da odrede mutagene aktivnosti i biodegradaciju ovih jedinjenja
- ekscitovana elektronska stanja i njihova uloga u fotodegradaciji polutanata
- nekovalentne interakcije

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Strani jezici engleski, nemački

Izabrani radovi Poglavlja

1. Bojana D. Ostojić, Dragana S. Đorđević: "The polycyclic aromatic hydrocarbons – fuel components and ubiquitous environmental pollutants" in Polycyclic aromatic hydrocarbons; Editor: Clifford Boone, Nova Science Publishers, 2015, ISBN: 978-1-63483-657-9
2. Bojana D. Ostojić, Dragana S. Đorđević: "The molecular properties of some diesel fuel components and their biodegradation" in Diesel Fuels: Characteristics, Performances and Environmental Impacts; Editors: Cristobal Silva and Agustin Rivera, Nova Science Publishers, 2013, ISBN: 978-1-62618-866-2
3. Bojana D. Ostojić, Dragana S. Đorđević: "The application of Density Functional Theory to calculation of properties of environmentally important species di- and trimethylnaphthalenes" in Density Functional Theory: Principles, Applications and Analysis, Editors: Joseph Morin and Jean Marie Pelletier, Nova Science Publishers, 2013, ISBN: 978-1-62417-954-9

Pregledni članak

1. Perić, M., Ostojić, B. and Radić-Perić, J. (1997) *Ab initio* investigation of the Renner-Teller effect in tetra-atomic molecules. Physics Reports: Review Section of Physics Letters 290, pp. 283-357

Radovi u časopisima

- Bojana D. Ostojić, Dragana S. Đorđević (2015) Electronic properties of environmental pollutants and their mutagenic activity: nitro derivatives of azaphenanthrenes, Chemosphere, Volume 135, pp. 319-324
- Bojana D. Ostojić, Dragana S. Đorđević (2015) Two nitro derivatives of azabenz[a]pyrene N-oxide: electronic properties and their relation to mutagenic activity, J. Hazardous Materials, Volume 285, pp. 94-102
- Bojana D. Ostojić, Branislav Stanković, Dragana S. Đorđević (2014) Theoretical study of the molecular properties of dimethylantracenes as properties for the prediction of their biodegradation and mutagenicity, Chemosphere, Volume 111, pp. 144-150
- Bojana D. Ostojić, Branislav Stanković, Dragana S. Đorđević (2014) The molecular properties of nitrobenzanthrone isomers and their mutagenic activities, Chemosphere, Volume 104, pp. 228-236
- Komainda, Adrian; Ostojic, Bojana; Koepfel, Horst (2013) *Ab Initio* Quantum Study of Nonadiabatic S1-S2 Photodynamics of S-Trans-Butadiene. J. Phys.Chem. A Volume 117, Issue 36, pp. 8782–8793
- Ostojic, Bojana; Jensen, Per; Schwerdtfeger, Peter; Bunker, Philip (2013): The Predicted Spectrum and Singlet-Triplet Interaction of the Hypermetallic Molecule SrOSr, J. Phys. Chem. A 117 (39), pp. 9370-9379
- Bojana D. Ostojić, Dragana S. Đorđević (2012) The electronic properties of trimethylnaphthalenes as properties for the prediction of biodegradation rates: *Ab initio* and DFT study, Chemosphere, 88(1), pp. 91-97
- Ostojić, B., Bunker, P. R., Schwerdtfeger, P., Assadollahzadeh, B. and Jensen, P. (2011) The predicted spectrum of the hypermetallic molecule MgOMg, Phys. Chem. Chem. Phys., 13, pp. 7546-7553
- Ostojić, B. D., Janjić, G. V., and Zarić, S. D. (2008) Parallel alignment of water and aryl rings – crystallographic and theoretical evidence for the interaction, Chem. Commun. 48, pp. 6546-6548
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- Bojana D. Ostojić, Slobodan Mišić, Dragana S. Đorđević (2013) A theoretical study of conformational flexibility, magnetic properties and polarizabilities of trimethylnaphthalenes, Int. J. Quantum Chem., vol. 113, pp. 1890-1898
- Bojana D. Ostojić, Dragana S. Đorđević (2012) *Ab initio* and density functional study of barrier heights for methyl group torsion and conformational deformability in 1,4,6-trimethylnaphthalene, Chem.Phys.Lett. 536, pp. 19-25
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- Bojana M. Francuski, Sladjana B. Novaković, Bojana D. Ostojić, Djordje D. Francuski, Goran A. Bogdanović (2015) Electronic features and hydrogen bonding capacity of the sulfur acceptor in thioureido-based

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- Bojana D. Ostojić, Branislav Stanković, and Dragana S. Đorđević (2014) Aromaticity and conformational deformability of some environmental pollutants - methylated anthracenes, *Fresenius Environmental Bulletin*, 23 (12) pp. 3036-3040
- B. Ostojić, P.R. Bunker, P. Schwerdtfeger, Artur Gertych, Per Jensen (2012) The predicted infrared spectrum of the hypermetallic molecule $CaOCa$ in its lowest two electronic states $X^1\Sigma_g^+$ and $a^3\Sigma_u^+$, *J. Mol.Struct.* 1023, pp. 101-107
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Saopštenja

1. B. Ostojić, P. R. Bunker, P. A. Schwerdtfeger, and P. Jensen, The predicted spectrum of the SbH_2 molecule in its two lowest Renner- degenerate electronic states $X^2B_1-A^2A_1$, The 24th Colloquium on High Resolution Molecular Spectroscopy HRMS, 2015, 24-28 August, 2015, Dijon, France.
2. Branislav Stanković, Bojana Ostojić, and Dragana Đorđević, The molecular properties of nitrodibenzofurans and their mutagenic activities, 18th International Symposium on Environmental Pollution and its Impact on Life in the Mediterranean Region; Crete, Greece, September 26 – 30, 2015.
3. B. Ostojić, Per Jensen, P. A. Schwerdtfeger, and P. R. Bunker, *Ab initio* calculations for the of group 2 M_2O hypermetallic oxides,, The 23rd International Conference on High Resolution Molecular Spectroscopy, Bologna, Italy, September 2-6, 2014
4. Branislav Stanković, Bojana Ostojić, and Dragana Đorđević, Theoretical investigation of molecular properties of methylanthracenes and biodegradation, 17th International Symposium on Environmental Pollution and its Impact on Life in the Mediterranean Region; Istanbul, Turkey, September 28 - Oct 1, 2013
5. Ostojić, B., Đorđević, D. The molecular properties of trimethylnaphthalenes and prediction of their biodegradation rates, 17th International Symposium on Environmental Pollution and its Impact on Life in the Mediterranean Region; Istanbul, Turkey, September 28 - Oct 1, 2013
6. P. Jensen, B. Ostojić, P. R. Bunker, P. Schwerdtfeger, and A. Gertych, P. Jensen, The predicted infrared spectrum of the hypermetallic molecule $CaOCa$ in its lowest two electronic states $X^1\Sigma_g^+$ and $a^3\Sigma_u^+$, The 22nd International Conference on High Resolution Molecular Spectroscopy, Prague, Czech Republic, 4-8 September, 2012.
7. B. Ostojić, P. R. Bunker, P. Schwerdtfeger, A. Gertych, and P. Jensen, The predicted infrared spectrum of the hypermetallic molecule $CaOCa$ in its lowest two electronic states $X^1\Sigma_g^+$ and $a^3\Sigma_u^+$, 67th International Symposium on Molecular Spectroscopy, Columbus, Ohio, USA, 18-20 June, 2012.
8. Ostojić, B., Jensen, P., Schwerdtfeger, P., Assodollahzadeh, B., and Bunker, P.R. The predicted infrared spectrum of the hyperberyllium molecule $BeOBe$ in its $X^1\Sigma_g^+$ and $a^3\Sigma_u^+$ electronic states, The 21st International Conference on High Resolution Molecular Spectroscopy, Poznan, Poland, 7-11 September 2010.
9. Yurchenko, S.D., Ostojic, B., Jensen, P., and Bunker, P.R., The predissociation of the d1A2 state of methylene: a combined *ab initio* (MRCI) and variational (TROVE), The 21st International Conference on High Resolution Molecular Spectroscopy, Poznan, Poland, 7-11 September 2010.
10. S. N. Yurchenko. P. Jensen, R.J. Buenker, P.R. Bunker, and B. Ostojić: The Near Ultraviolet Band System of Singlet Methylene, The 17th Colloquium on High Resolution Molecular Spectroscopy, Papendal, The Netherlands, 9-13 September 2001.