



**51. SAVETOVANJE  
SRPSKOG HEMIJSKOG DRUŠTVA  
i  
2. KONFERENCIJA MLADIH HEMIČARA SRBIJE**

**PROGRAM**

*i*

**KRATKI IZVODI RADOVA**

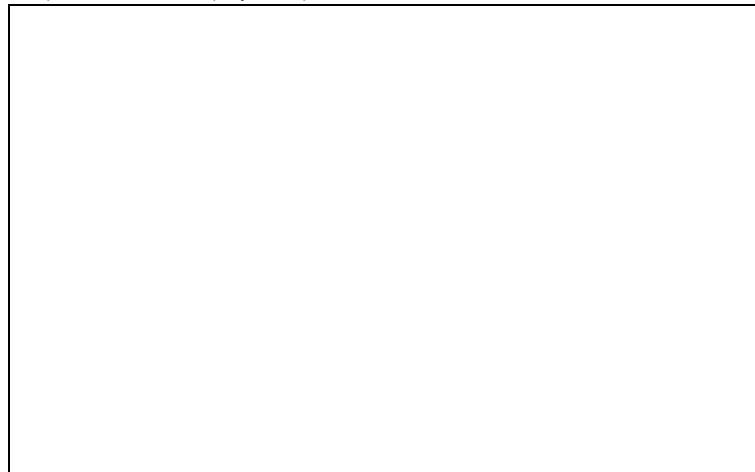
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and  
2<sup>nd</sup> Conference of the Young Chemists of Serbia**

**Program**

**&**

**Book of Abstracts**

*Niš, 5-7. juni 2014.  
Niš, Serbia, June 5-7, 2014*



**51. SAVETOVANJE SRPSKOG HEMIJSKOG DRUŠTVA i  
2. KONFERENCIJA MLADIH HEMIČARA SRBIJE, NIŠ, 5-7. JUNI 2014.  
PROGRAM I KRATKI IZVODI RADOVA**

5<sup>ST</sup> MEETING OF THE SERBIAN CHEMICAL SOCIETY i  
2<sup>ND</sup> CONFERENCE OF THE YOUNG CHEMISTS OF SERBIA, NIŠ, SERBIA, JUNE 5-7, 2014  
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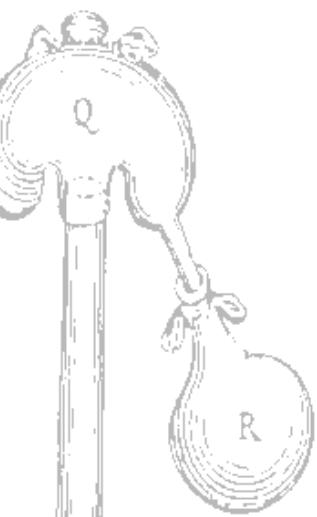
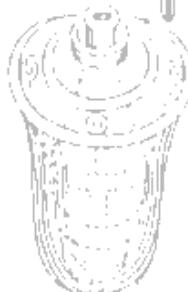
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# **51. SAVETOVANJE SRPSKOG HEMIJSKOG DRUŠTVA**

**51<sup>st</sup> Meeting of the Serbian Chemical Society**

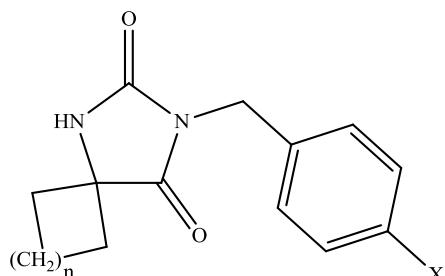
## **KRATKI IZVODI ABSTRACTS**

**Sinteza, struktura i svojstva novih  
3-(supstituisanih benzil)-5,5'-cikloalkan spirohidantoina**

Anita M. Lazić, Bojan Đ. Božić, Nemanja P. Trišović, Nataša V. Valentić,  
Gordana S. Uščumlić

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U radu je sintetizovano dvadeset novih 3-(supstituisanih benzil)-5,5'-cikloalkan-spirohidantoina čija je struktura određena temperaturama topanja, FT-IR,  $^1\text{H}$  i  $^{13}\text{C}$  NMR i UV spektroskopijom. Solvatochromizam ovih jedinjenja je proučavan metodom linearne korelacije energija solvatacije. Kvantitativna procena uticaja rastvarača na pomeranje apsorbacionih maksimuma je nedvosmisleno ukazala na vezu između lipofilnosti i relevantnih farmakoloških karakteristika proučavanih jedinjenja sa njihovim proton-akceptorskim svojstvima.



$n = 2, 3, 4$

$X = \text{H, CH}_3, \text{OCH}_3, \text{Cl, Br, CN, NO}_2$

*Slika. Struktura proučavanih jedinjenja  
Figure. Structure of the investigated compounds*

**Synthesis, structure and properties of novel  
3-(substituted benzyl)-5,5'-cycloalkanespirohydantoins**

Anita M. Lazić, Bojan Đ. Božić, Nemanja P. Trišović, Nataša V. Valentić,  
Gordana S. Uščumlić

*Faculty of Technology and Metallurgy, University of Belgrade, Karnegijeva 4, Belgrade, Serbia*

In this work twenty novel 3-(substituted benzyl)-5,5'-cycloalkanespirohydantoins have been synthesized and fully characterized by melting point, FT-IR,  $^1\text{H}$  and  $^{13}\text{C}$  NMR and UV spectroscopy. Solvatochromism of these compounds have been analysed by means of linear solvation energy relationships (LSER) concept. The quantitative relationships between hydrogen bonding interaction and the corresponding pharmacologically relevant properties of the studied compounds have been discussed.