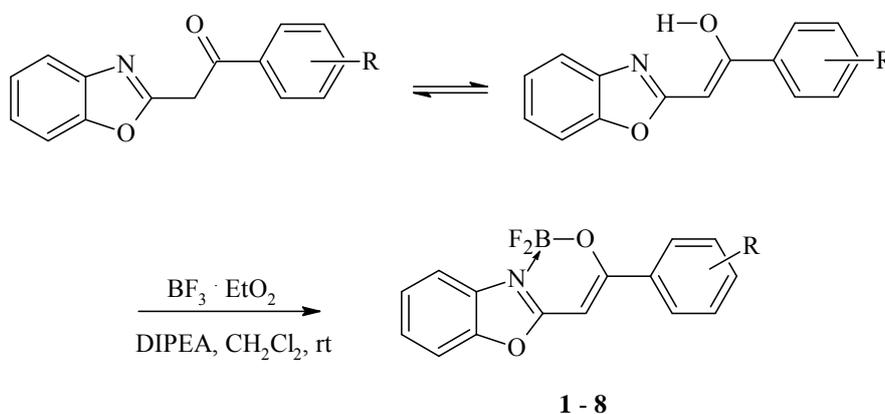


# SYNTHESIS, STRUCTURE AND PROPERTIES OF THE DIFLUOROBORANES OF SUBSTITUTED 2-PHENACYLBENZOXAZOLES AS POTENTIAL PHOTOSENSITIZERS IN THE PHOTOPOLYMERIZATION PROCESS

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Organoboron complexes are one of the most important types of fluorescent dyes. The proton involved in intramolecular hydrogen bonding of substituted 2-phenacylbenzoxazoles can be easily replaced by another acid such as the  $\text{BF}_2^+$  cation. The presented study provides a report on the basic photophysical properties of a recently synthesized series of substituted 2-phenacylbenzoxazole difluoroboranes (Scheme 1) with a view towards the effect of substitution on their properties [1].



**Scheme 1.** The schematic representation of the synthesis of 2-phenacylbenzoxazole difluoroboranes where R = 4-NMe<sub>2</sub> (**1**), 4-OMe (**2**), 4-Me (**3**), 3-Me (**4**), H (**5**), 3-OMe (**6**), 4-Cl (**7**), 3-Cl (**8**).

Complexes have been identified based on a magnetic atomic nucleus <sup>1</sup>H, <sup>11</sup>B, <sup>13</sup>C, <sup>15</sup>N, and <sup>19</sup>F isotope resonance spectra. The photophysical properties of these new dyes have been investigated by fluorescence and ultraviolet-visible (UV-Vis) absorption spectroscopy. The changes in Stokes shifts and the molar absorption coefficients related to the nature of the substituent in 2-phenacyl moiety are also reflected in changes in the fluorescence quantum yields. The study using these compounds as photosensitizers in the photopolymerization process is still in progress.