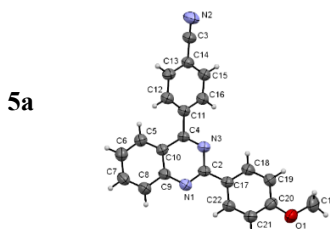
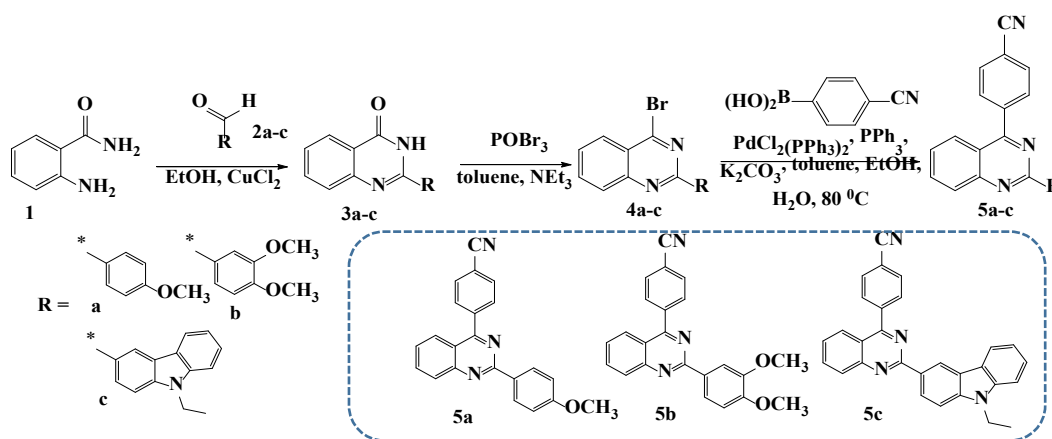


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SYNTHESIS AND PHOTOPHYSICAL PROPERTIES
OF 2-ARYL-4-(4-CYANOPHENYL)QUINAZOLINES

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Abstract. The establishment of detailed structure–property relationships (SPRs) provides beneficial information for fine-tuning of the key characteristics and the rational design and synthesis of fluorophores. Quinazoline derivatives containing electron donating fragment at position 4 of quinazoline core were described previously by our research group.[1,2,3] Some quinazolines demonstrated strong fluorescent properties in solution as well as solvatochromic and sensing properties.

In this work we have designed and synthesized 4-(4-cyanophenyl)-quinazoline counterparts with electron donating unit at position 2. The target products **5a-c** were obtained by Pd-catalyzed cross coupling reaction between 4-bromo-derivatives **4a-c** and 4-cyanophenylboronic acid.



	Solvent	λ_{abs} , nm	λ_{em} , nm	QY, %
5a	MeCN	348, 292, 258	505	<0.1
	Toluene	352, 297	447	<0.1
5b	MeCN	314, 280	595	< 0.1
	Toluene	360, 319	500	10.9
5c	MeCN	300	326, 421	< 0.1
	Toluene	379	500	23.2

The compounds were purified by column chromatography and the structure was confirmed by NMR-spectroscopy, mass-spectrometry and X-ray analysis data. The photophysical properties were measured in MeCN and toluene solutions, the fluorescent quantum yield is up to 23%.

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