

Theoretical modelling of the frequency-dependent phase shift between emission and excitation in ZnO nanorod photoluminescence spectra

Radim Novák¹, Zdeněk Remeš¹, Štěpán Remeš¹, and Júlia Mičová²

¹*Institute of Physics of the Czech Academy of Sciences, Na Slovance 1999/2, 182 00 Praha 8, Czechia*

²*Institute of Chemistry SAS, Dubravska cesta 9, 84538 Bratislava, Slovakia*

Hydrothermally growth, zinc oxide (ZnO) nanocrystals are known for their interesting scintillation and photocatalytic properties. In our previous paper we presented an optical setup for measuring spectrally resolved photoluminescence (PL) mean decay time using conventional UV LED with sinusoidal excitation and a phase shift method with about 10 ns time resolution [1], recently upgraded to 1 ns. Photoluminescence decay is the process by which the excited state of a material, induced by the absorption of photons, returns to the ground state, emitting light in the process. This process can be measured in time domain, with the decay rate indicating how quickly the material returns to its ground state after excitation, or equivalently in frequency domain, with the difference between the phase of the light wave at the moment of excitation and the phase of the light wave at the moment of emission. Here we report on the theoretical modelling of the frequency-dependent phase shift between emission and excitation in photoluminescence spectra of ZnO nanorods.

This work was supported by the CAS Open Science 2024, CAS-SAS Mobility Plus project SAV-23-13, CAS-SAS-2022-08, MSTC Danube project 8X23025 and by the CSF project 24-10607J. This work was also supported by the Slovak Research and Development Agency under the Contract no. DS-FR-22-0035.

[1] Z Remeš, M Buryi, J Pejchal, O Babčenko, Š Remeš, R Novák and J Mičová, J. Phys.: Conf. Ser. 2712 (2024) 012004