

Study of Perfluorophosphonic Acid Surface Modifications on Zinc Oxide Nanoparticles

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Table S1. Binding energies determined by XPS.

	ZnO control	PFPDPA	5FBPA	F₂₁DDPA
	Binding energy (eV)	Binding energy (eV)	Binding energy (eV)	Binding energy (eV)
Zn2p_{3/2}	1022.0	1021.9	1021.9	1021.9
Zn2p_{1/2}	1045.1	1045.0	1045.0	1044.9
O1s	530.6	530.6	530.4	530.6
	531.8	531.9	532.0	531.7
C1s		284.8	287.9	293.3
		285.5	285.6	291.1
		287.7		288.4
				285.5
P2p		133.6	134.1	134.1
F1s		688.0	688.1	688.2
		685.0	685.0	684.7

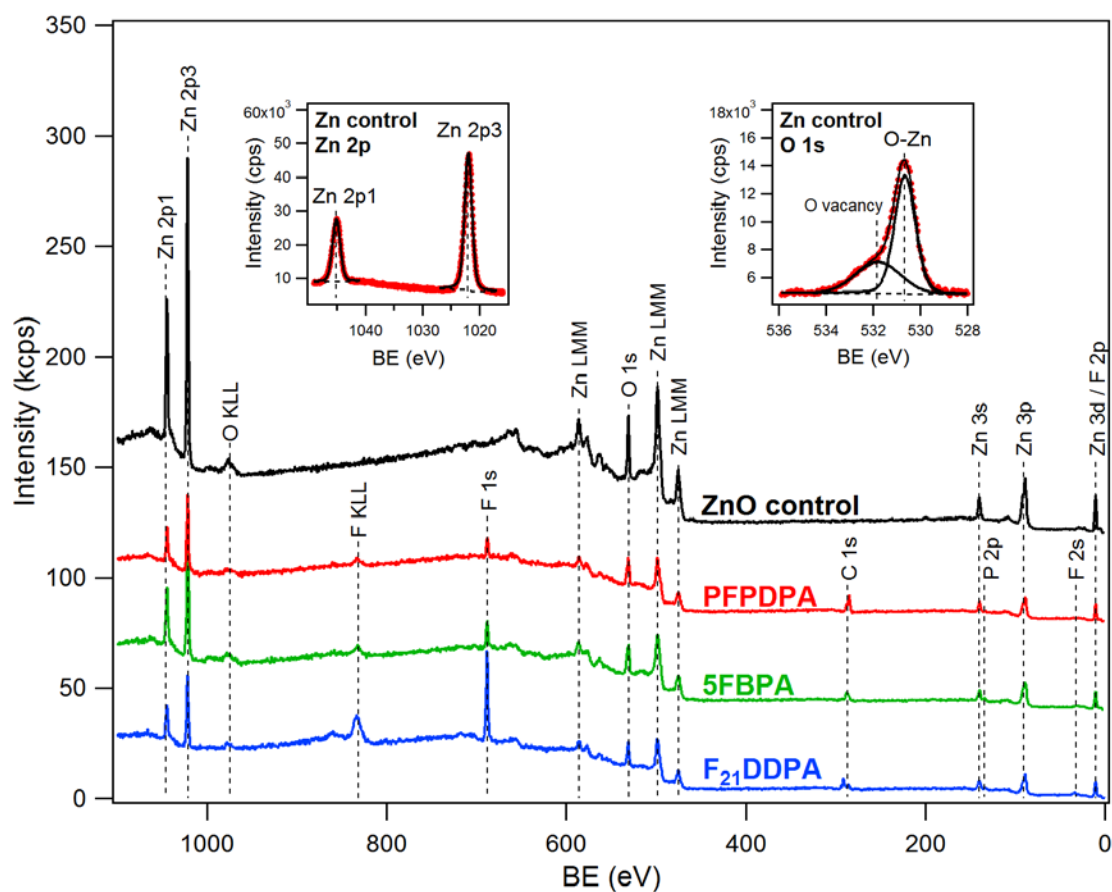


Figure S1. XPS survey spectra of (a) ZnO nanoparticles unmodified (control) (black line), (b) PFPDPA on ZnO (red line), (c) 5FBPA on ZnO (green line), and (d) F₂₁DDPA on ZnO nanoparticles (blue line). High-resolution core level spectra of ZnO unmodified (black) for Zn2p and O1s are also shown.

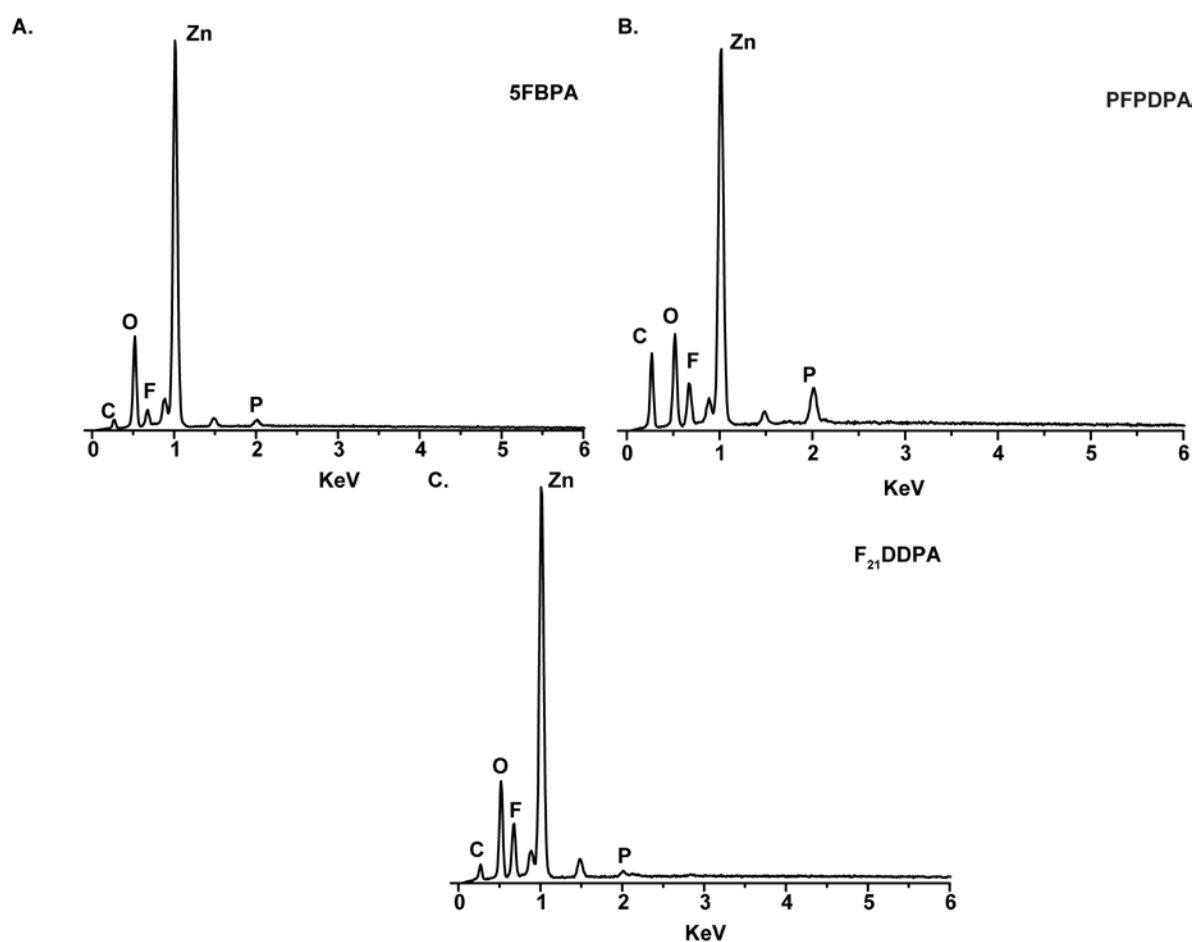


Figure S2. EDS spectra for (a) 5FBPA, (b) PFPDPA, and (c) F₂₁DDPA modified ZnO nanoparticles.