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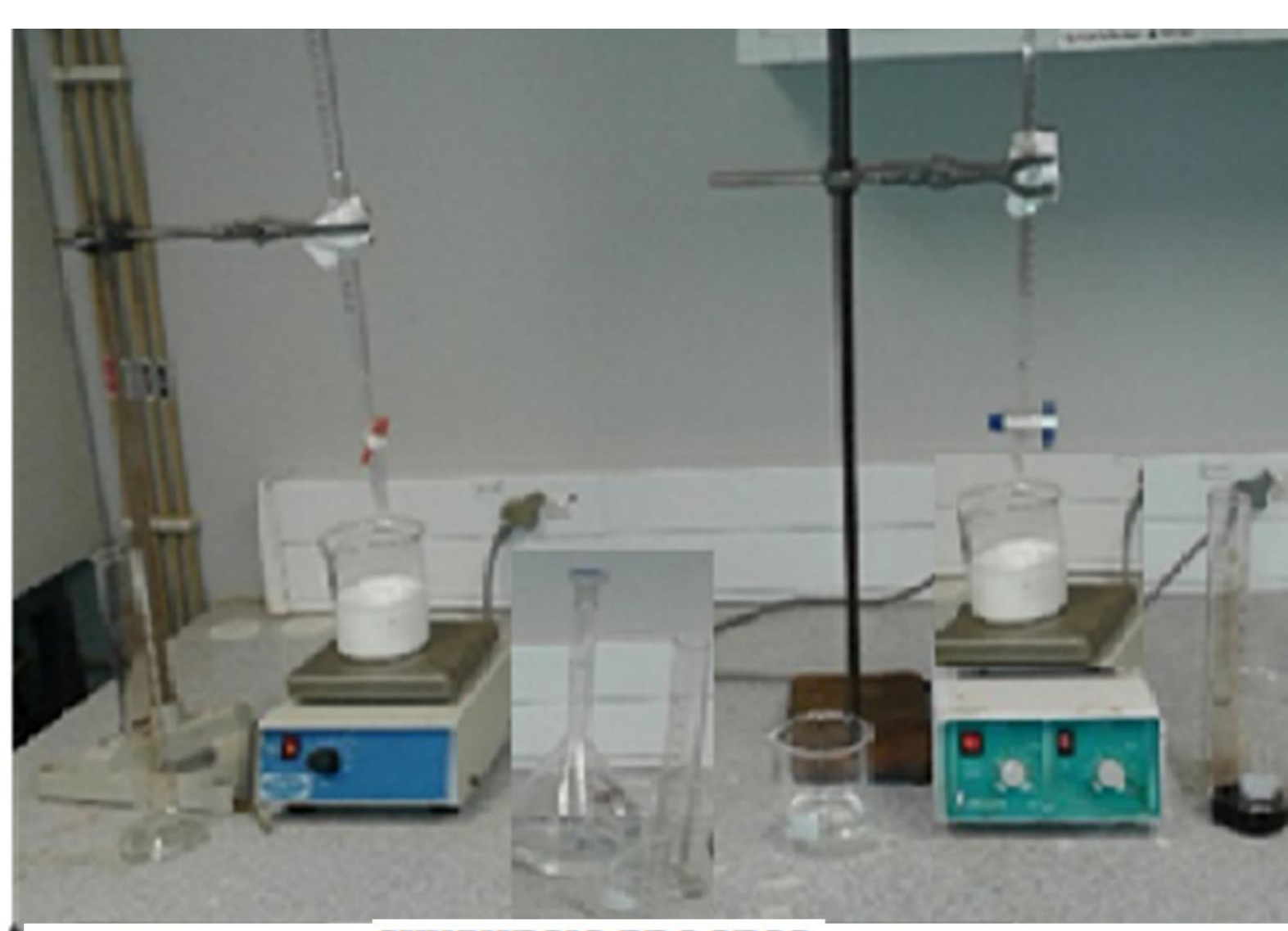
CdMgZnO NANOPARTICLES CHARACTERIZATION AND RESULTS ANALYSIS

INTRODUCTION

The red shift of the basic absorption peak centered at 375 nm (3.30 eV) to 432 nm (2.92 eV) in the UV-visible absorption spectra validates band gap narrowing of the CdMgZnO of nanoparticles. Microstructural experiment and structural adjustment with lower angle of diffraction peaks and vibrational stretching of Zn-O respectively in the lattice from the substitution of Cd makes it a suitable and a possible host for nanophosphor application in solid state lighting [1]

Furthermore, the impact of divalent metals (Mg and Cd) on semiconductor oxide nanostructure with respect to luminescence emissions in the visible spectra has the potential of being applied in light emitting devices, laser devices, non-linear optical devices, flat panel displays, sensors, anti-reflecting coatings and IR windows [2]

EXPERIMENTAL PROCESS



SYNTHESIS PROCESS

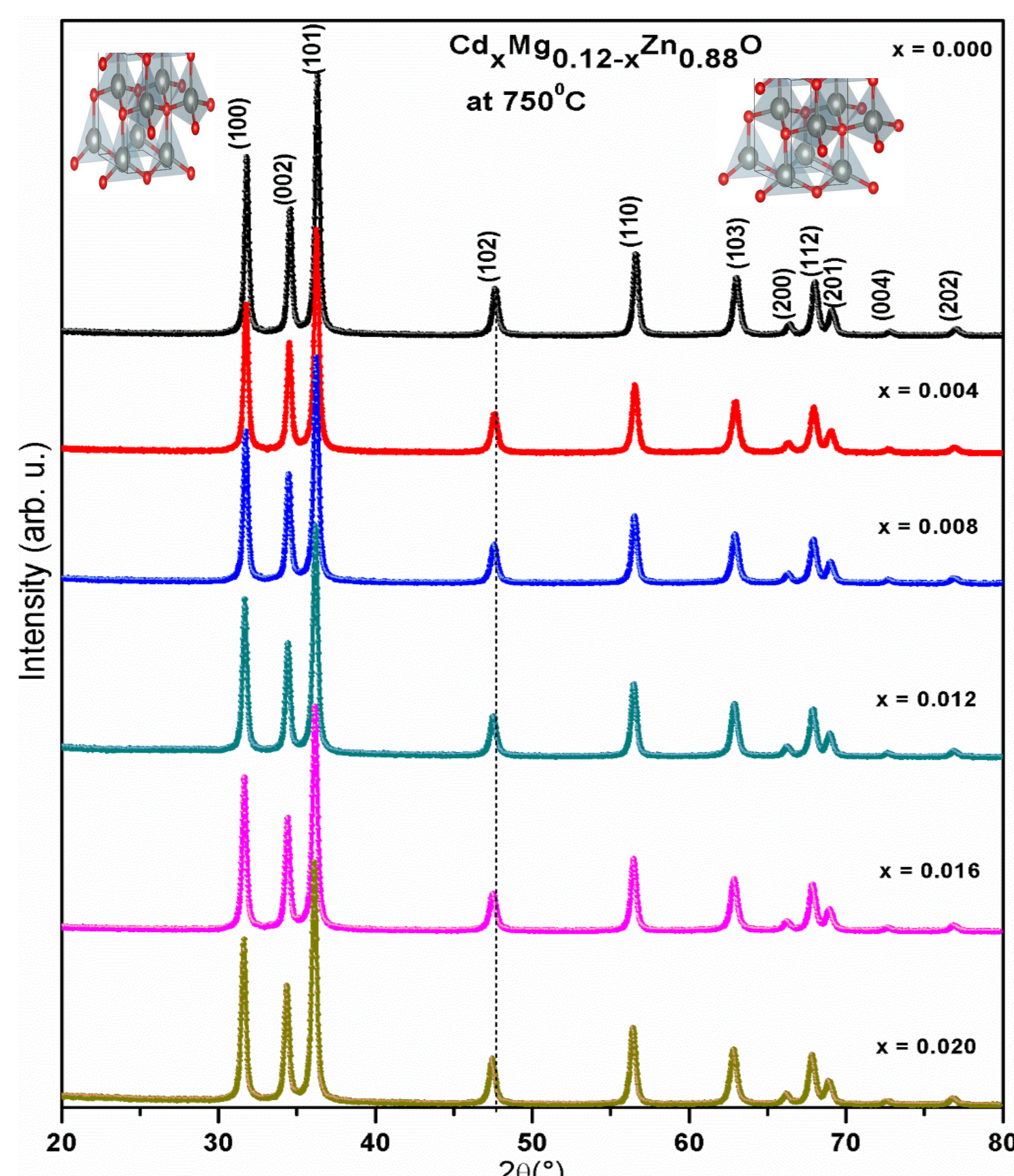
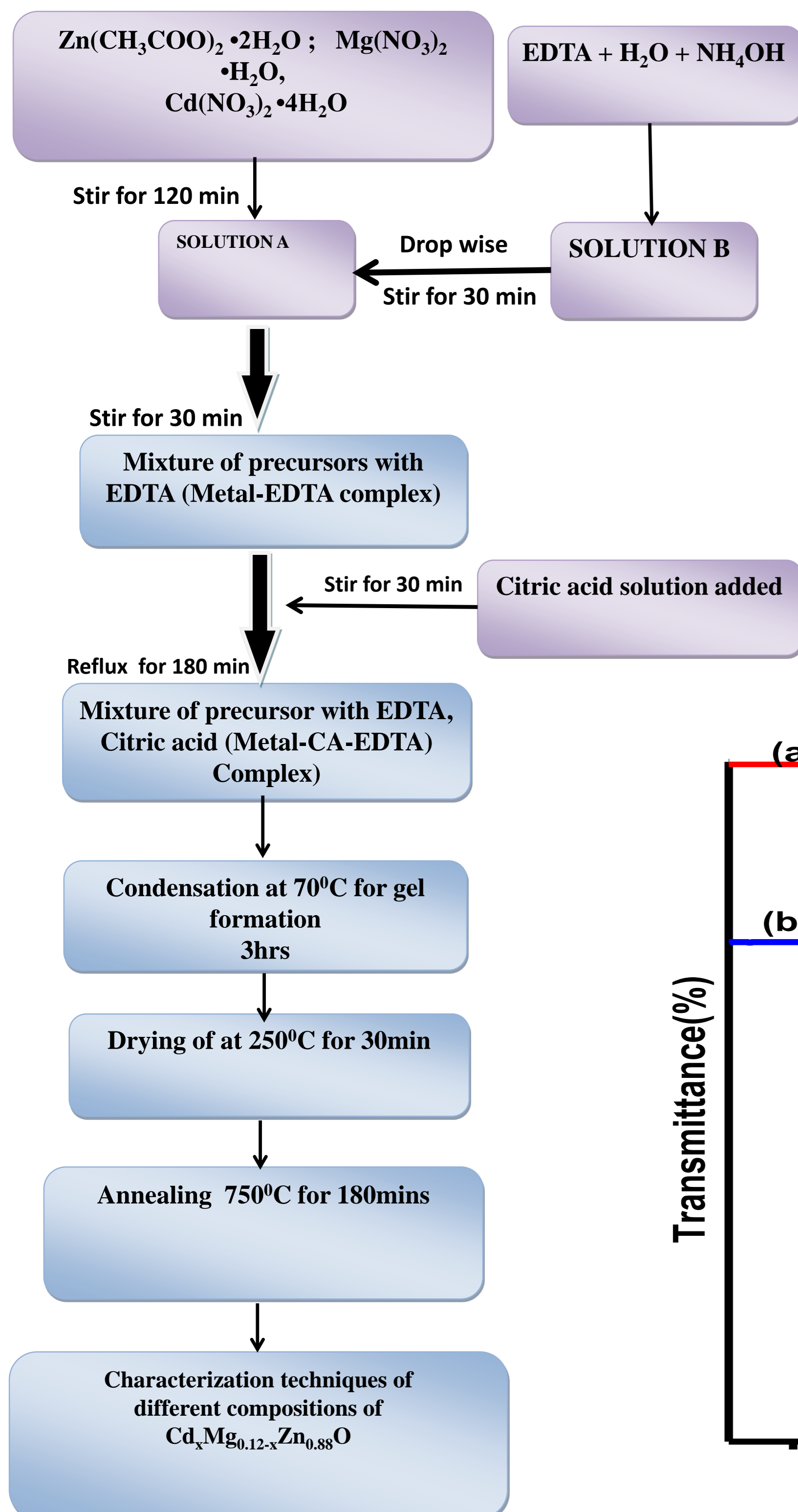


Figure 1: Plots of XRD Pattern of CdMgZnO.

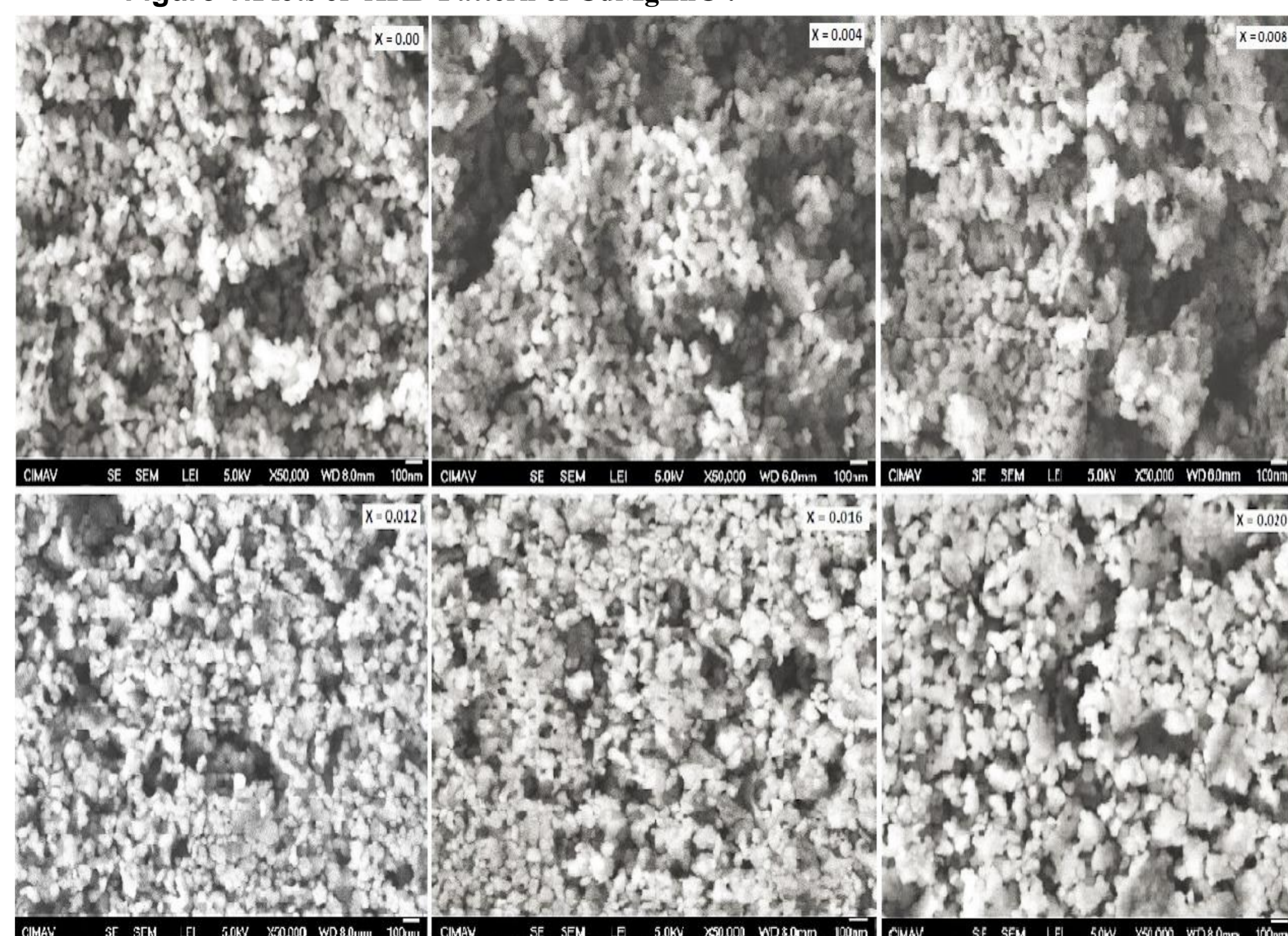


Figure 3: (a) SEM micrographs of CdMgZnO nanoparticles

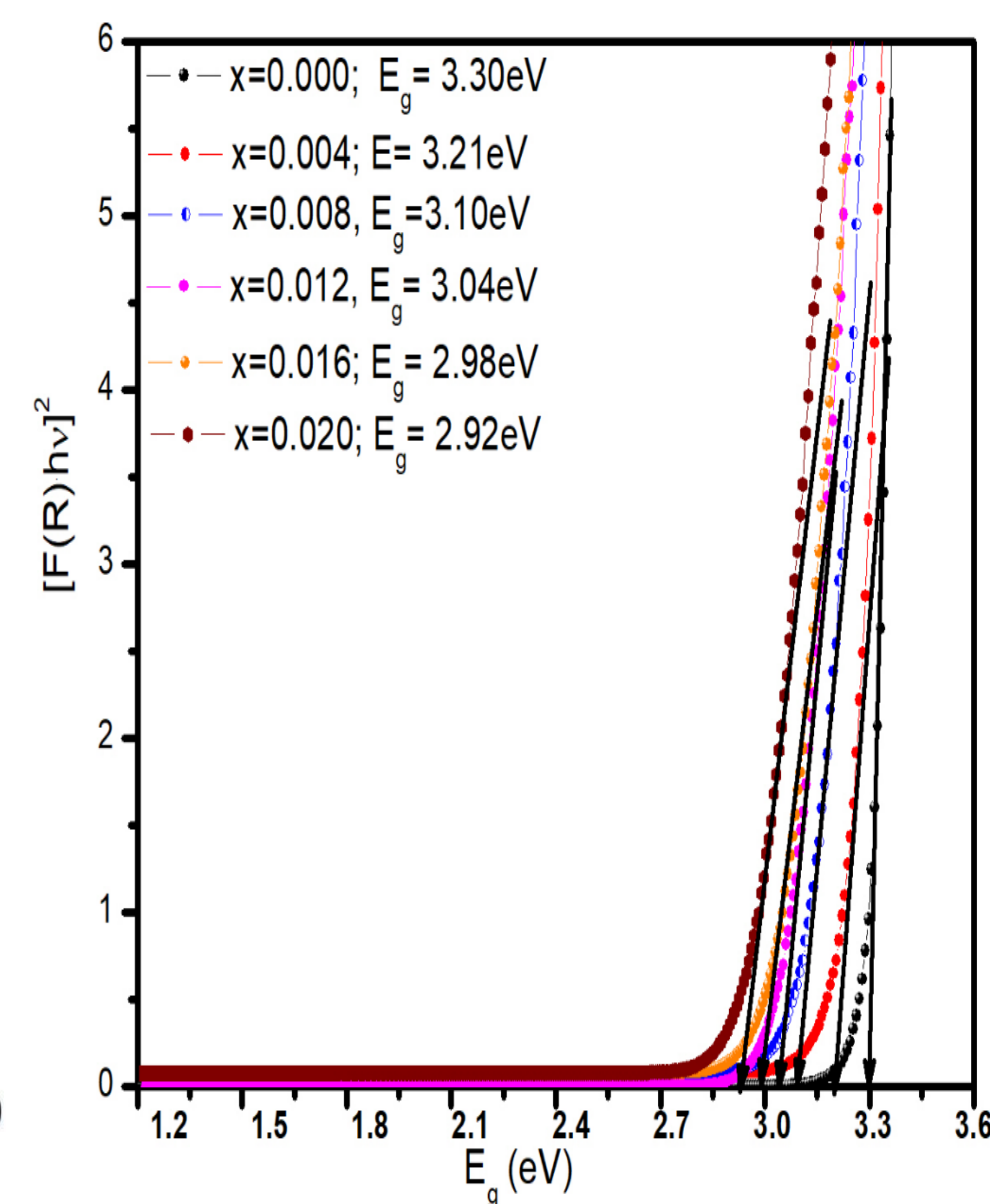
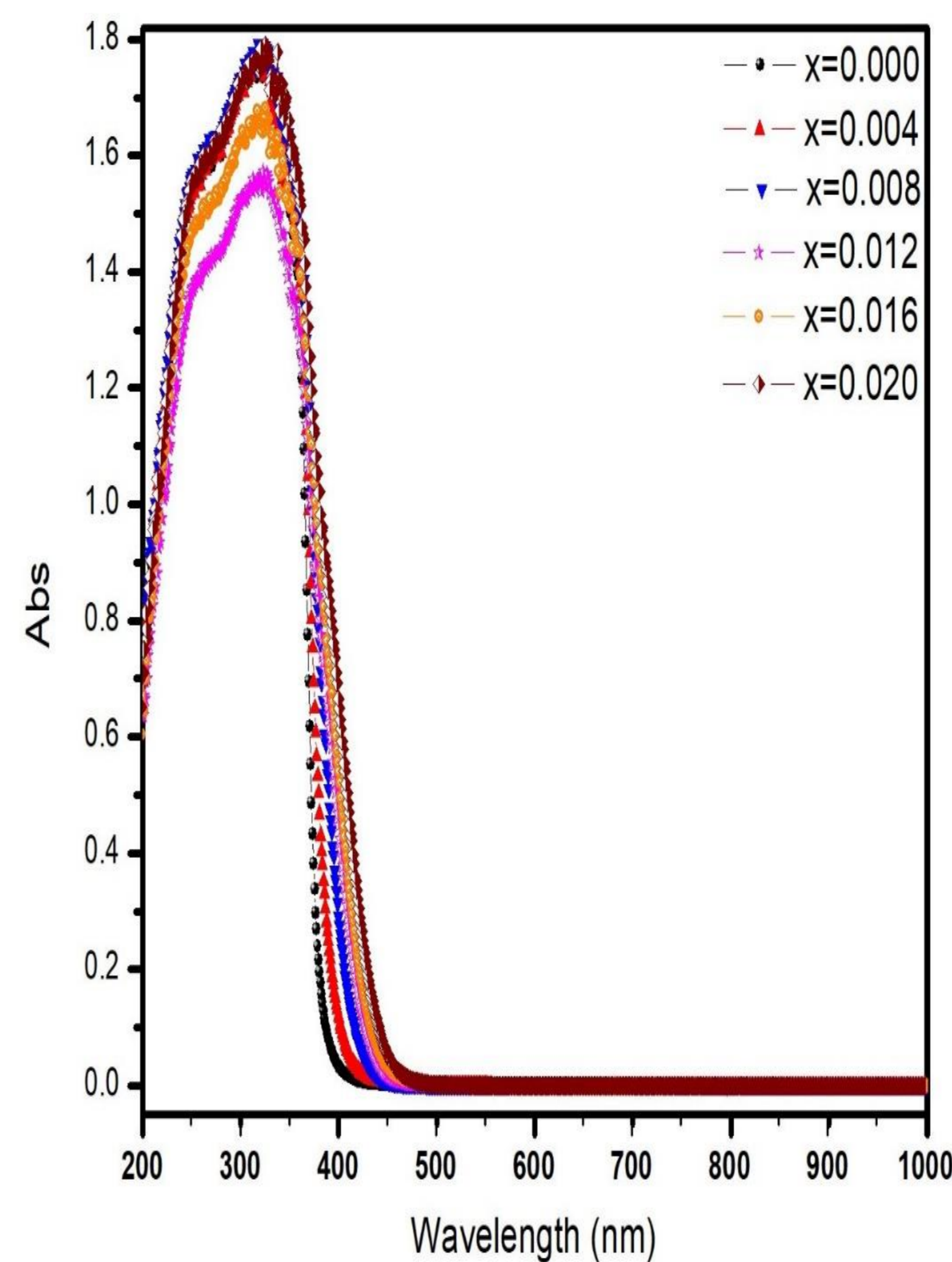


Figure 2: UV-Vis-NIR absorbance and optical band-gap of CdMgZnO nanoparticles

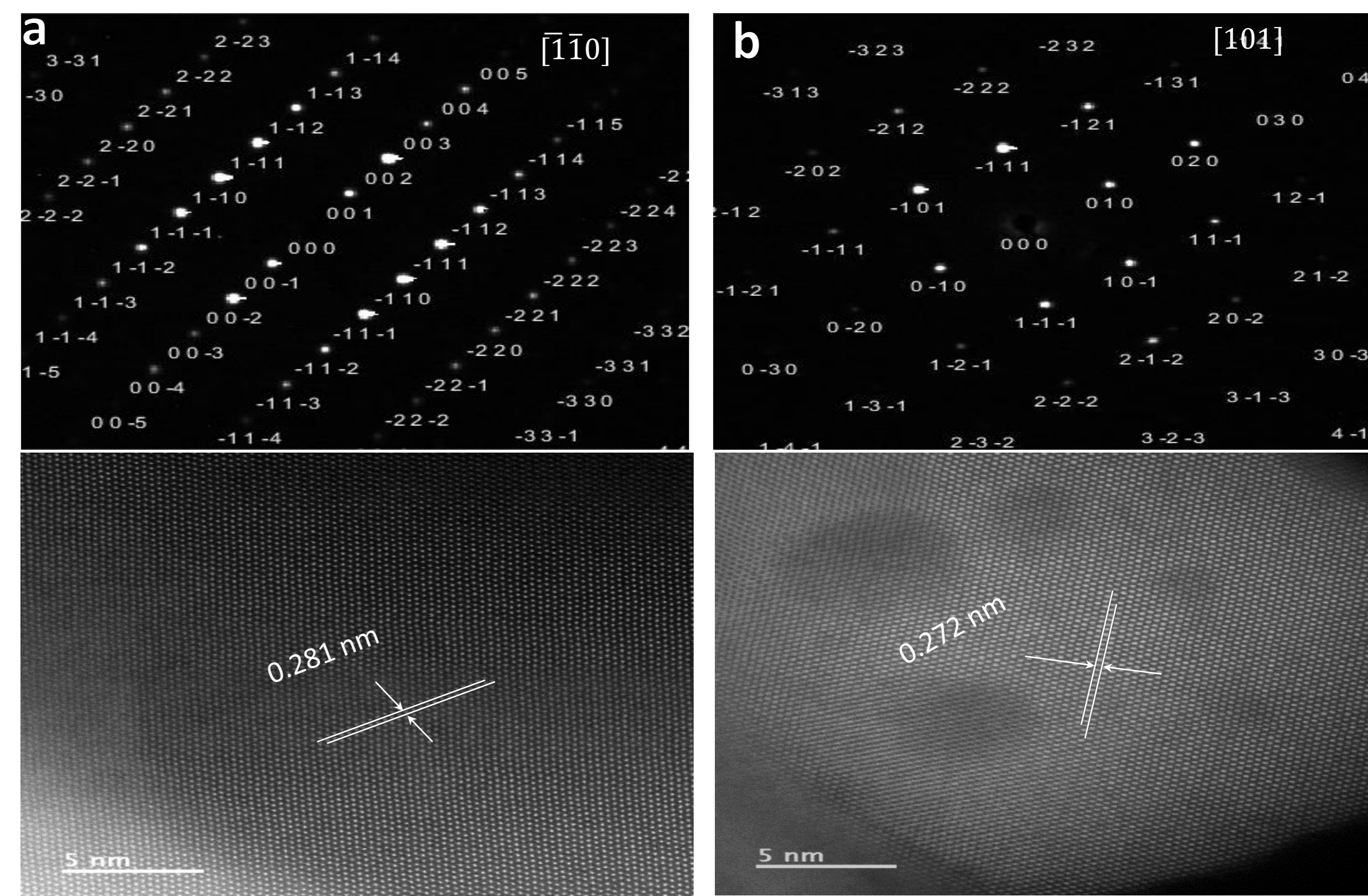


Figure 4: TEM images of CdMg_(0.12-x)Zn_{0.88}O nanoparticles. a) SAED pattern in the zone axes $[1\bar{1}0]$ and HRTEM for $x = 0.02$ (maximum concentration). b) SAED pattern in the zone axes $[101]$ and HRTEM for $x = 0$ (without Cd concentration).

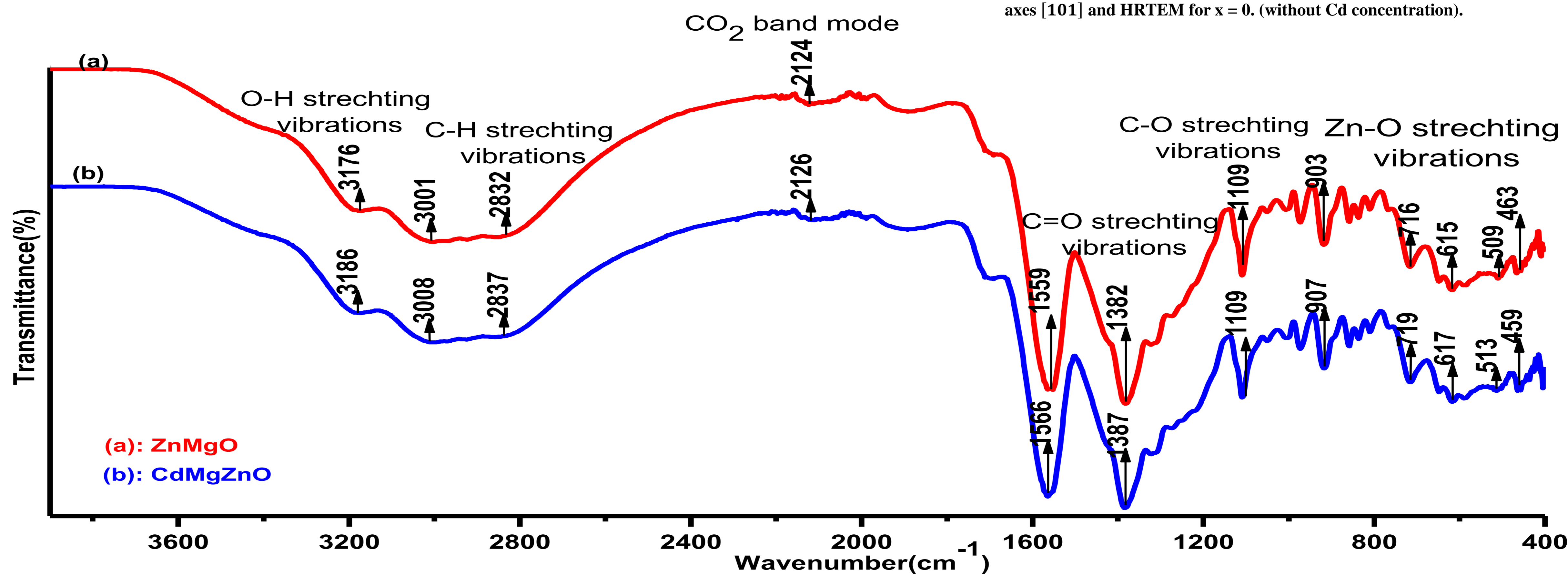


Figure 5: FTIR of Cd_xMg_(0.12-x)Zn_{0.88}O nanoparticles ($x = 0.000, 0.020$).

DISCUSSION

The entire XRD pattern matched to a (JCPDS 36-1451) standard, revealed a hexagonal wurtzite structure with a single phase crystal [2] as shown in Fig.1(a) with a degree of stress or strain in the synthesized nanoparticles since Cd²⁺ ion has a larger ionic radius of R_{Cd}^{2+} (0.095nm) to that of R_{Zn}^{2+} (0.074nm) and R_{Mg}^{2+} (0.072nm) respectively.

The increase in the absorption wavelength may be linked with the presence of impure states below the conduction band from the Cd concentration with increase in crystallite sizes from the XRD.

Furthermore, repulsion effect between Cd 4d and O 2p levels in the valence band is a consequence of narrowing the optical band gap

FTIR absorption peak around 465 cm⁻¹ assigned to Zn-O bond stretched at 513 cm⁻¹ and 903 cm⁻¹ due to the effect of Mg and Cd content with narrower peaks for CdMgZnO nanoparticles

CONCLUSION

In this study the dependent of features devices on the properties of nanomaterials with the substitution of divalent metals (Cd and Mg) into ZnO is an effective route for improving and modifying the optical band gap and can be considered as a material for near UV and blue optoelectronic application which is confirmed from the decreased in optical band gap which varied from 3.30 eV to 2.92 eV with particle size increase and preservation of the crystal symmetry of the hexagonal wurtzite structure.

REFERENCES
 [1] O. Kalu, J. A. D. Moller, A. R. Rojas, Structural and optical properties of cadmium magnesium zinc oxide (CdMgZnO) nanoparticles synthesized by sol-gel method, Phys. Lett. A, 383(10) (2019) 1037-1046. <https://doi.org/10.1016/j.physleta.2018.11.0>
 [2] L.T. Jule, F.B. Dejene, A.G. Ali, K.T. Roro, A. Hegazy, N. K. Allam, E.E. Shenawy Wide visible emission and narrowing band gap in Cd-doped ZnO nanopowders synthesized via sol-gel route, J. Alloys compds. 687 (2016) 920-926.

PUBLISHED ARTICLE
 [1] O. Kalu, J. A. D. Moller, A. R. Rojas, Structural and optical properties of cadmium magnesium zinc oxide (CdMgZnO) nanoparticles synthesized by sol-gel method, Phys. Lett. A, 383(10) (2019) 1037-1046. <https://doi.org/10.1016/j.physleta.2018.11.0>

Table 1. Analysis of crystallographic parameter of Cd_xMg_(0.12-x)Zn_{0.88}O (0.000 ≤ x ≤ 0.020) nanoparticles, crystallites sizes (Scherrer's and W-H method) and strain

Con. of Cd ²⁺ (x) at. %	d spacing (Å)			Lattice parameters			Cell volume (V)Å ³	Crystal size (nm)		Strain (ε × 10 ⁻³)
	(100)	(002)	(101)	a = b (Å)	c (Å)	c/a		Scherrer's	W - H	
0.00	2.8078	2.5913	2.4688	3.2421	5.1827	1.5985	47.1792	20.06	20.68	1.60
0.004	2.8079	2.5921	2.4698	3.2422	5.1841	1.5989	47.1907	20.15	21.06	1.44
0.008	2.8082	2.5929	2.4700	3.2425	5.1859	1.5993	47.2163	20.20	21.56	1.12
0.012	2.8116	2.5967	2.4723	3.2461	5.1939	1.6000	47.3680	20.46	22.61	1.01
0.016	2.8127	2.5993	2.4732	3.2474	5.1976	1.6005	47.4679	20.65	23.81	0.83
0.020	2.8128	2.5999	2.4738	3.2479	5.1984	1.6005	47.4889	21.11	26.15	0.53