Supplementary Information (SI) for Journal of Materials Chemistry C. This journal is © The Royal Society of Chemistry 2024

Supplementary File



S1: Elemental study using EDS spectroscopy:

Figure S1: (a-d) EDS spectra with corresponding elemental composition of Z0, Z1, Z2 and Z3 samples, respectively.

S2: UV-Visible study-

The room temperature optical absorption (UV-Vis) spectra of GO, rGO and CZFO/rGO (Z0, Z1,Z2 and Z3) nanocomposites are shown in Figure S2. In Figure S2 the optical absorption spectra of GO and rGO shows a sharp absorption peak around 221nm. This peak is related to the π - π * transitions for aromatics -C-C- originating for sp² domains [1]. For pure Co_{0.8}Zn_{0.2}Fe₂O₄ MNPs in figure 9(b) significant absorption shows in the visible range. After adding of rGO in Co_{0.8}Zn_{0.2}Fe₂O₄ MNPs in different wt %, the main absorption peak was observed at around 200-230nm ,which can be attributed

to the intrinsic absorption of rGO and $Co_{0.8}Zn_{0.2}Fe_2O_4$ MNPs respectively .The optical energy band (E_g) of all the sample compositions was calculated from the absorption spectra by using the Tauc plot method follows; $(\alpha hv)^n = A'$ (hv-E_g), where hv indicates the energy of the incident photon, A' is a constant and n shows an index that is characterizes of the optical absorption process and its value is n = 2 and 1/2 based on the direct and indirect band gap transitions, respectively [2]. Here, the value of n taken $\frac{1}{2}$ for direct band gap transitions. This depends on the nature of the electronic transmission responsible for the reflection in the materials. The absorption coefficient α can be calculated by the Beer Lambert Law [3, 4].

I=I_o e^{- α t}, where $\alpha = 2.303 \times A/t$

Where A is the absorbance and t are thickness (1cm) of standard quartz cuvette which we have used during experiment. The energy band of pure GO, rGO and $Co_{0.8}Zn_{0.2}Fe_2O_4/rGO$ (Z0, Z1, Z2 and Z3) nanocomposites have been calculated to be about 2.7, 2.8,2.5,2.3,2.1 and 2.6 eV respectively. The calculated energy band of pure rGO and $Co_{0.8}Zn_{0.2}Fe_2O_4/rGO$ nanocomposites corresponds to the result to previous study. On the other hand pure $Co_{0.8}Zn_{0.2}Fe_2O_4$ MNPs have reduced the energy band gap of $Co_{0.8}Zn_{0.2}Fe_2O_4/rGO$ nanocomposites as the increases the rGO % content ,the energy band gap change from 2.5 to 2.1 eV and further increases 2.6 eV. This reduction of energy band gap of $Co_{0.8}Zn_{0.2}Fe_2O_4/rGO$ nanocomposites with different rGO doping concentration is due to the increase of photoexcited electron-hole in the $Co_{0.8}Zn_{0.2}Fe_2O_4/rGO$ nanocomposites with different rGO and composites with increases the rGO nanocomposites with increases the rGO and smaller crystallites size of MNPs [32].



Figure S2: (a) UV-Vis absorption spectra (b) energy band gap of pure GO, rGO and CZFO/rGO (Z1,Z2 and Z3) nanocomposites at room temperature.

Reference:

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