## Comment on "Molten Salt Synthesis of Bi<sub>2</sub>WO<sub>6</sub> Powders and its Visible-Light Photocatalytic Activity" in Materials Research. 2019; 22(5): e20190311

## Salmon Landi Jr.ª 回

<sup>a</sup>Instituto Federal de Educação Ciência e Tecnologia Goiano, 75901-970, Rio Verde, GO, Brasil.

Received: April 22, 2021; Accepted: June 24, 2021

Recently, Dai et al. published a paper in Research Materials 2019; 22(5): e20190311. They have investigated the visible-light photocatalytic activity of the synthesized  $Bi_2WO_6$  powders<sup>1</sup>. In the context of determining of band gap energy based on diffuse reflectance measurements, the cited authors make some mistakes, which are clarified in the present letter.

Keywords: Diffuse Reflectance Spectroscopy, Kubelka-Munk function, Absorption coefficient.

- 1)  $\alpha hv = A(hv E_g)^2$  was designed by Kubelka-Munk (K-M) equation<sup>1</sup>, which is incorrect. In fact,  $\alpha hv = A(hv - E_g)^2$  is a classical equation from the optical transitions theory in semiconductor/insulator materials<sup>2</sup>. On the other hand, the dimensionless K-M function/equation (*F*(*R*)) is defined as the ratio between the K-M absorption and scattering coefficients<sup>3</sup>. In the vicinity of  $E_g$ , the scattering phenomenon can be neglected, that is, the scattering coefficient is treated as a constant<sup>4</sup>, therefore: *F*(*R*)  $\propto \alpha$
- According to the referred authors<sup>1</sup>, the parameter *A* is usually 1. In fact, the value of *A* does not interfere in the determination of the band gap energy (*E<sub>g</sub>*) because it is an independent parameter of photon energy (*hv*)<sup>5</sup>. However, its value depends on properties such as the number of unit cell or even the type of electronic transition between the valence and conduction bands<sup>6</sup>. In this sense, it is unlikely that its value is usually 1.
- 3) According to inset of Figure 5,  $(\alpha hv)^{1/2}$  is plotted against the hv, where  $\alpha$  is the absorption coefficient of material and represents the probability of light being absorbed per unit path length<sup>7</sup>. Consequently, the unit of measure for  $\alpha hv$  is unit of energy divided by unit of length and not eV as declared by the authors<sup>1</sup>. However, as the values presented on the ordinate axis in the inset shown in Figure 5 were obtained from the absorbance measurements, the most appropriated unit for  $\alpha hv$  is "arbitrary unit".

## References

- Dai B, Xuan M, Lv Y, Jin C, Ran S. Molten Salt Synthesis of Bi<sub>2</sub>WO<sub>6</sub> Powders and its Visible-Light Photocatalytic Activity. Mater Res. 2019;22(5):e20190311. http://dx.doi. org/10.1590/1980-5373-mr-2019-0311.
- Landi S. Comment on "Photocatalytic degradation of RhB from an aqueous solution using Ag3PO4/N-TiO2 heterostructure" and "Evaluation of the effect of dose change of Fe3O4 nanoparticles on electrochemical biosensor compatibility using hydrogels as an experimental living organism model". J Mol Liq. 2021;338:116635. http://dx.doi.org/10.1016/j. molliq.2021.116635.
- Milosevic M, Berets SL. A review of FT-IR of diffuse reflection sampling considerations. Appl Spectrosc Rev. 2002;37(4):347-64. http://dx.doi.org/10.1081/ASR-120016081.
- Liu S, Wang Z, Bao Q, Li X, Chen Y, Wang Z, et al. Abnormal thermal quenching and blue-shift of Zn3(BO3) (PO4): Inducing host T defect by doping Mn<sup>2+</sup> and Tb<sup>3+</sup>. Dyes Pigments. 2021;165:44-52. http://dx.doi.org/10.1016/j. dyepig.2019.01.048.
- Malainho E, Vasilevskiy MI, Alpuim P, Filonovich SA. Dielectric function of hydrogenated amorphous silicon near the optical absorption edge. J Appl Phys. 2009;106:073110. http://dx.doi. org/10.1063/1.3240203.
- Elliott RJ. Intensity of Optical Absorption by Excitons. Phys Rev. 1957;108:1384-9. http://dx.doi.org/10.1103/PhysRev.108.1384.
- Yang L, Kruse B. Revised Kubelka-Munk theory. I. Theory and application. J Opt Soc Am A Opt Image Sci Vis. 2004;21:1933-41. http://dx.doi.org/10.1364/josaa.21.001933.