Erratum: Electronic Structures and Optical Properties of Fe_2VAl and Fe_2VGa Alloys

[J. Korean Phys. Soc. **47**, 692 (2005)]

Joo Yull RHEE*

BK21 Physics Research Division and Institute of Basic Science, Sungkyunkwan University, Suwon 440-746

DOI: 10.3938/jkps.62.1708

Figure 3 should be swapped by the following figure.

Fig. 3. Theoretical optical conductivity spectra of Fe_3Al and Fe_2VAl alloys. The λ -fitting is not applied. The OC spectrum of Fe_2VGa alloy is also included for comparison.

*E-mail: rheejy@skku.edu