



Determination of Mass Attenuation Coefficients, Effective atomic number and Electron Density of Lumefantrine in the Energy Range 1 keV – 100 GeV

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Abstract

The effective atomic number and electron density of Lumefantrine(LU) have been calculated for total and partial photon interactions by the direct method in the wide energy range of 1 keV – 100 GeV using WinXCOM. The values of these parameters have been found to change with energy. The variations of effective atomic number and electron density with energy are calculated and shown graphically.

Keywords: LU; Mass attenuation coefficients; Effective atomic number; Electron density.

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