MOSSBAUER STUDY OF Fee. oiNio. 99Cr2S4

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Abstract - Feo. or Nio. 99 Cr2S4 has been studied by Mössbauer spectroscopy and X-ray diffraction. The crystal structure is found to be monoclinic with the lattice parameters: a = 5.910 Å, b = 3.410 Å, c = 11.11 Å and β = 91.50 °. Magnetic hyperfine and quadrupole interactions in the antiferromagnetic state at 81 K have been studied, yielding the following results: H = 156 kOe, 1/2eqQ(1+1/3 η ²) $^{1/2}$ = -2.15 mm/s, θ = 59 °, ϕ = 90 °, and η = 0.9. Electronic energy levels of Feo.orNio.99Cr2S4 can be explained in terms of a Hamiltonian involving crystal field energy, spin-orbit couplings, and exchange interactions.