
Adsorption of N₂, O₂, CO₂ and H₂ on Hydrotalcite-Like System: Mg²⁺-Al³⁺-(Fe(CN)₆)⁴⁻

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Abstract: The compound Mg²⁺-Al³⁺-(Fe(CN)₆)⁴⁻, with a structure similar to hydrotalcite Mg₆Al₂(OH)₁₆CO₃·4H₂O, is prepared by a coprecipitation method. Chemical, thermal, and X-ray analysis of the compound lead to the formula: (Mg_{0.741}Al_{0.259}(OH)₂)(Fe^{II}(CN)₆)_{0.067}(CO₃)_{0.002}(H₂O)_{1.05})

The compound is dehydrated at 100° to 250° C and the adsorption of N₂, O₂, CO₂, and H₂ on it is measured at -196° C to room temperature by the volumetric method. The water is removed reversibly from 100° to 200° C and the interstices after dehydration act as sites of adsorption. When dehydration is carried out at 150° C the adsorptive activity reaches a maximum. Adsorption isotherms of N₂ and O₂ at -196° C and of CO₂ at room temperature are of the Langmuir type, and the saturated amounts of N₂, O₂, and CO₂ adsorbed are 96.3, 65.2, and 91.8 ml (STP)/g, respectively.

Adsorption isotherms of N₂ and O₂ at room temperature are of the Henry type. The amount of O₂ adsorbed is about 67% of that of N₂. The isosteric heats of adsorption at room temperature of N₂, O₂, and CO₂ are 5.1, 4.8, and 11.0 kcal/mol, respectively. A dehydrated product of a Mg²⁺-Al²⁺-CO₃²⁻ compound does not permit adsorption of CO₂ at room temperature, but permits the adsorption of H₂O. Hence, it has a molecular sieve effect. No adsorption of H₂ is observed in any of the compounds tested.

Key Words: Adsorption • Hydrotalcite • Isotherm • Surface • Zeolite

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