

H₂, O₂和CO在有机分子和聚合物上吸附的CNDO/2研究

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摘要 应用CNDO/2法研究了H₂, O₂和CO在吡咯和由吡咯单元构成的模型大分子上的吸附行为,讨论了被吸附分子相对于杂环烃的不同吸附取向的影响,优化了各个体系相对于能量的分子间距,考察了相互作用能随分子间距和吡咯基单元的数目,得到了一系列与相关实验相互印证的有用的规律。

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CNDO/2 study on the adsorption behavior of hydrogen, oxygen and carbon monoxide on organic molecules and polymers

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Abstract The CNDO/2 method was used to study the adsorption behavior of H₂, O₂, and CO on pyrrole and model macromolecules built up from pyrrole units. Different orientations of the adsorption of the adsorbed molecule relative to the heterocyclic hydrocarbon were considered. The intermolecular distance was optimized with respect to the energy for each composite system. The dependence of the interaction energy on the intermolecular distance and on the number of pyrrole units was examined. A series of useful rules in agreement with experimental findings were obtained.

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