

硅沸石完美骨架上的吸附 II. 乙胺、正戊烷、乙醇在 FAU型沸石上的吸附热

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摘要 用测定吸附等温线法研究乙胺、正戊烷和乙醇在疏水高硅FAU沸石上的吸附热效应。根据Clapeyron-Clausius方程, 处理吸附等温线, 得到不同覆盖度 C (C =被吸附分子数/晶胞)的等量吸附热 $Q_{st}(C)$ (一定覆盖度 C 时由Clapeyron-Clausius方程计算的吸附热)及平均等量吸附热 Q_{st}^* (一定温度区间里等量吸附热 Q_{st} 的平均值), 以及 Q_{st}^* 与沸点蒸发热 ΔH_v 的差值 ΔH_1 (定义为相互作用强度 $\Delta H_1 = Q_{st}^* - \Delta H_v$)。所研究的三种有机分子的 ΔH_1 的次序为 $\Delta H_1(\text{乙胺}) > \Delta H_1(\text{正戊烷}) > \Delta H_1(\text{乙醇})$ 。这与AT值(定义为脱附温度 T_d 与吸附质的沸点温度 T_b 的差值, 无需单位)有正相关关系。由AT值观察到的高硅FAU沸石Si-O骨架与被吸附乙胺之间可能存在的强相互作用, 为本研究测定的热力学定量数据 ΔH_1 值所证明。

关键词 [吸附](#) [等量吸附热](#) [乙胺](#) [戊烷](#) [乙醇](#) [高硅沸石](#) [吸附等温线](#) [吸附热](#)

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Adsorption on siliceous zeolites with perfect framework II. adsorption heat of ethylamine, n-pentane and ethanol on FAU zeolite

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Abstract The adsorption heat of ethylamine, n-pentane and ethanol was measured based upon the determination of adsorption isotherms at different temperatures on high silica FAU type zeolite with perfect Si-O framework. Calculated with Clapeyron - Clausius equation, the isosteric heat of adsorption $Q_{st}(C)$ (adsorption heat at certain coverage, which is defined as the numbers of adsorbed molecules per unit cell) and average isosteric heat of adsorption Q_{st}^* (the average adsorption heat with various coverage determined in a certain temperature range) were obtained for the adsorbates investigated. ΔH_1 -the difference of Q_{st}^* and ΔH_v (the heat of vaporization of certain adsorbate), was defined as interaction strength for characterizing the interaction between Si-O framework of the zeolite and the adsorbed organic compound. The order of ΔH_1 for the adsorbed organic compounds is as follows: $\Delta H_1(\text{ethylamine}) > \Delta H_1(\text{n-pentane}) > \Delta H_1(\text{ethanol})$, which is consistent with the order of AT value (affinity index, was defined as $AT = T_d - T_b$, where T_d is the temperature of the weight loss peak of the DTG curve, and T_b is the boiling point of a certain adsorbate). The strong interaction between Si-O framework and the adsorbed ethylamine, which was observed from the desorption behavior on FAU zeolite, is thermodynamically proved in this study by the high interaction strength of $\Delta H_1(\text{ethylamine})$ determined quantitatively.

Key words [ADSORPTION](#) [ETHYLAMINE](#) [PENTANE](#) [ETHANOL](#) [HIGH SILICA ZEOLITES](#) [ADSORPTION ISOTHERM](#) [ADSORPTION HEAT](#)

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