CXN天然沸石的研究2:吸附性质

李军,邱瑾,龙英才

复旦大学化学系.上海(200433)

收稿日期 修回日期 网络版发布日期 接受日期

摘要 采用N~2,NH~3,CO~2,乙烯,丙烯,水,甲醇,乙醇,丙醇等作为吸附剂,研究了由我国CXN天然沸石改性制得的H-STI和Na-STI沸石的吸附性质,H-STI和Na-STI沸石的BET表面积及微孔孔体积约为420m^2/g和0.20m^3/g。根据NH~3和CO~2在H-

STI沸石上的吸附等温线计算得到它们的吸附热分别为44.8和26.5kJ/mol。乙烯, 丙烯, 甲醇, 乙醇, 丙醇等在Na-

STI沸石上的吸附等温线表明该沸石对有机分子的吸附具有链长选择性。在低分压下水相对于甲醇的吸附量表明沸石具有一定的疏水性质。 关键词 沸石 吸附 选择性 氦 氦 二氧化碳 乙烯 丙烯 水 甲醇 乙醇 丙醇

分类号 0612

## Studies on CXN natural zeolite 2: Adsorption properties

Li Jun,Qiu Jin,Long Yingcai

Fudan Univ, Dept Chem.Shanghai(200433)

Abstract The adsorption properties of H-STI and Na-STI, which were prepared from the natural CXN zeolite obtained in China, were investigated with N $\sim$ 2, NH $\sim$ 3, CO $\sim$ 2, ethylene, propylene, water, methanol, ethanol and 1-propanol. The BET surface area and pore volume of H-STI and Na-STI were about 420 m $^{\circ}$ 2/g and 0.20 m $^{\circ}$ 3/g. The adsorption heat of NH $\sim$ 3 and CO $\sim$ 2 on H-STI calculated by the adsorption isotherms at different temperature was 44.8 and 26.5 kJ/mol respectively. The investigation of the adsorption isotherms of ethylene, propylene, methanol, ethanol and 1-propanol on Na-STI reveals the unusual selectivity to the length of the carbon chain in the molecules of organics. The loading of water is lower than that of methanol in lower partial pressure, which exhibits a somewhat hydrophobic property of the zeolite.

Key words ZEOLITE ADSORPTION SELECTIVITY NITROGEN AMMONIA CARBON DIOXIDE ETHYLENE PROPENE WATER METHANOL ETHANOL PROPANOL

DOI:

通讯作者

## 扩展功能

## 本文信息

- ► Supporting info
- ▶ <u>PDF</u>(0KB)
- ▶[HTML全文](0KB)
- ▶参考文献

# 服务与反馈

- ▶把本文推荐给朋友
- ▶加入我的书架
- ▶加入引用管理器
- ▶复制索引
- ► Email Alert
- ▶文章反馈
- ▶ 浏览反馈信息

#### 相关信息

▶ 本刊中 包含"沸石"的 相关文章

## ▶本文作者相关文章

- 李军
- 邱瑾
- · <u>龙英</u>才