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CO₂ Adsorption Performance of Functionalized Metal-Organic Frameworks of Varying Topologies by Molecular Simulations

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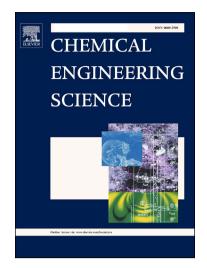
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CO₂ Adsorption Performance of Functionalized

Metal-Organic Frameworks of Varying Topologies

by Molecular Simulations

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KEYWORDS Carbon capture; adsorption; topology; pore size; functionalization

ABSTRACT

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