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Investigation of calixarene complexes with dasatinib and lapatinib: A Density Functional Theory approach

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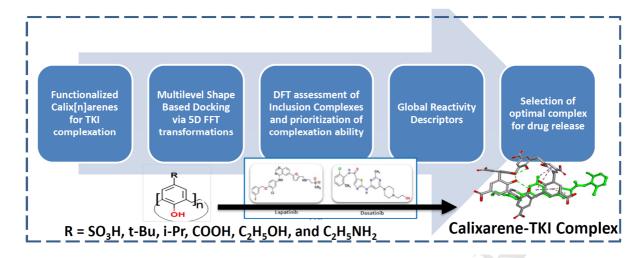
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