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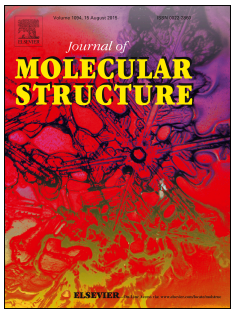

Molecular dynamics simulation of functionalized graphene surface for high efficient loading of doxorubicin

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