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Theoretical study of the substituent effect on the hydrogen atom transfer mechanism of *meta*- and *para*-substituted benzenetellurols

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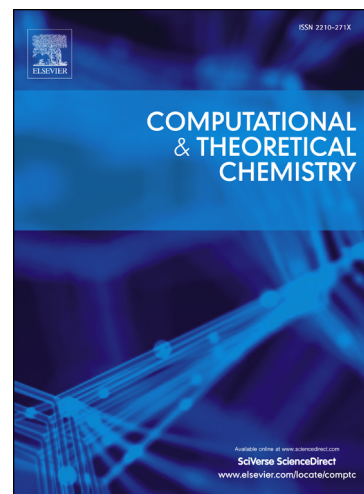
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