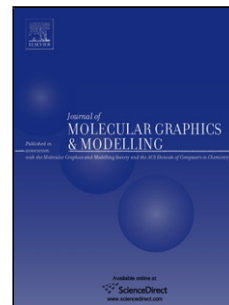


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ab initio molecular orbital calculations in water**

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