

Structural properties of Iridium (III) in water by ab initio quantum mechanical charge field molecular dynamics simulations

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Abstract	
Authors	P Iswanto, R Armunanto, HD Pranowo
Journal Name	The 1 st International Conference on Computation for Science and Technology, 75, 2010
Publish Year	2010
Citation	(not set)
Url	<a (iii)="" ab="" by="" charge="" dynamics="" field="" href="https://scholar.google.com/scholar?q=+intitle:" in="" initio="" iridium="" mechanical="" molecular="" of="" properties="" quantum="" simulations"="" structural="" water="">https://scholar.google.com/scholar?q=+intitle:"Structural properties of Iridium (III) in water by ab initio quantum mechanical charge field molecular dynamics simulations"
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