

Calculations of the Potential Energy Surface for a Water Molecule Dissociation on the Pt (111) Surface

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Abstract	
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Journal Name	Soedirman International Conference on Mathematics and Applied Sciences, 2022
Publish Year	2022
Citation	(not set)
Url	<a (111)="" a="" calculations="" dissociation="" energy="" for="" href="https://scholar.google.com/scholar?q=+intitle:" molecule="" of="" on="" potential="" pt="" surface="" surface"="" the="" water="">https://scholar.google.com/scholar?q=+intitle:"Calculations of the Potential Energy Surface for a Water Molecule Dissociation on the Pt (111) Surface"
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