

GolP-CHARMM: First-Principles Based Force Fields for Au(111) and Au(100)

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Citation Report

#	ARTICLE	IF	CITATIONS
9	A molecular dynamics simulation study of the electric double layer and capacitance of [BMIM][PF6] and [BMIM][BF4] room temperature ionic liquids near charged surfaces. <i>Physical Chemistry Chemical Physics</i> , 2013, 15, 14234.	1.3	93
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