



Volume 19, Issue 3

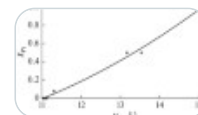
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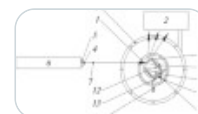
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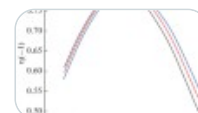
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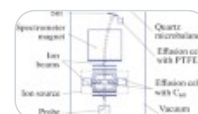
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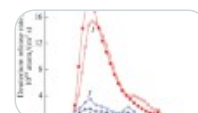
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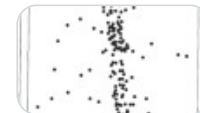


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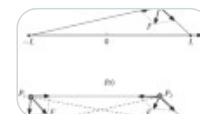
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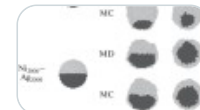
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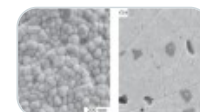
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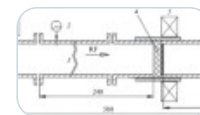
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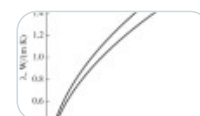
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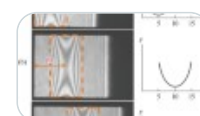
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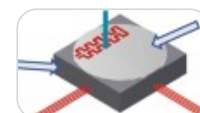
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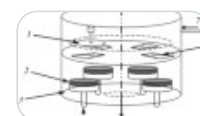
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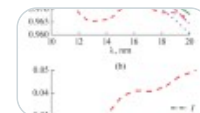
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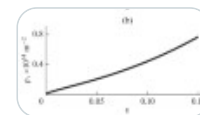
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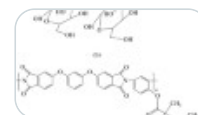


Monte Carlo Modeling of the Graphene Moiré Structure on an Ir(111) Substrate



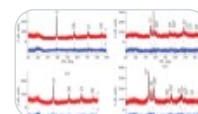
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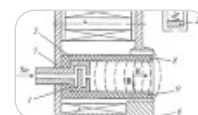
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Hydride Phases Based on High-Entropy Alloys $\text{TiZrTaV}(\text{Mo}_{1-x}\text{Nb}_x)$ and $\text{TiZrTaNb}(\text{Mo}_{1-x}\text{Ni}_x)$, $x = 0.2, 0.4, 0.6$, and 0.8



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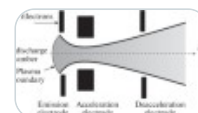
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