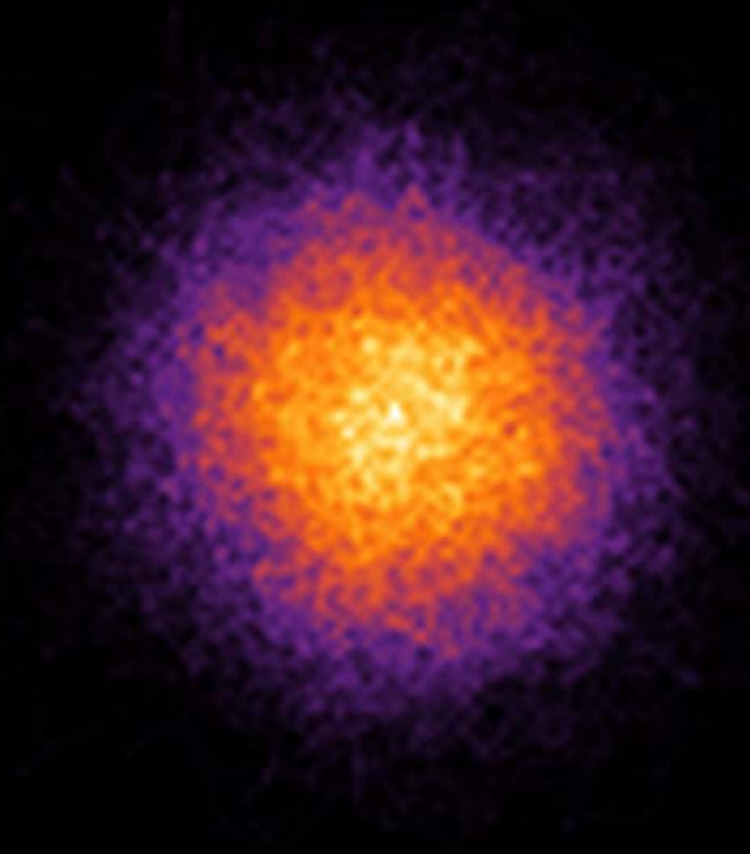


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# Development of molecular dynamics methodology for simulations of hard materials

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