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Molecular dynamics simulations of high energy cascade in ordered alloys: defect production and subcascade division

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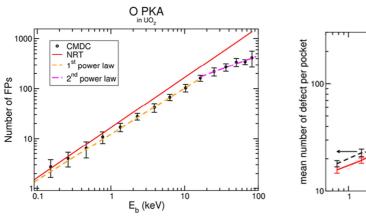
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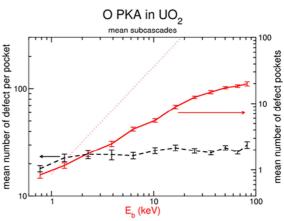


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Molecular Dynamics simulations of cascade in ordered alloys show that defect production in such materials does not follow the rules which apply for monoelemental solids

The equivalence between linearity of defect production with ballistic energy and apparition of subcascades does not hold in general for alloys.





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