

Abstract Submitted
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Coulomb Explosion of Transition Metal Oxides¹ D.E. BLUMLING,
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tosecond pulses of light (624 nm) are employed to investigate the formation of high
charge states and kinetic energy release (KER) from the Coulomb explosion of var-
ious transition metal oxide clusters. A molecular dynamics simulation is used to
predict KER values for the ground state structures of the representative species as
determined by density functional theory calculations.

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