Generalized Koopmans density functional calculations reveal the deep acceptor state of N_O in ZnO

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$E\,N$

f 2.1

II. METHOD

f. 1

$$V\,s = \lambda\,s$$

IV. HYBRID-FUNCTIONAL RESULTS FOR N_O

$\varepsilon_A(0/-)=E_V+1.$

$\Delta \quad \underset{s}{\equiv} +0. \quad 2 \quad . \quad^2 \quad s, \quad \quad f \quad s \quad \quad \quad s$

$\alpha = 0.25.$

f 1.

V

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