Supplementary Material

Density functional theory study of triple transition metal cluster anchored on the C_2N monolayer for nitrogen reduction reaction

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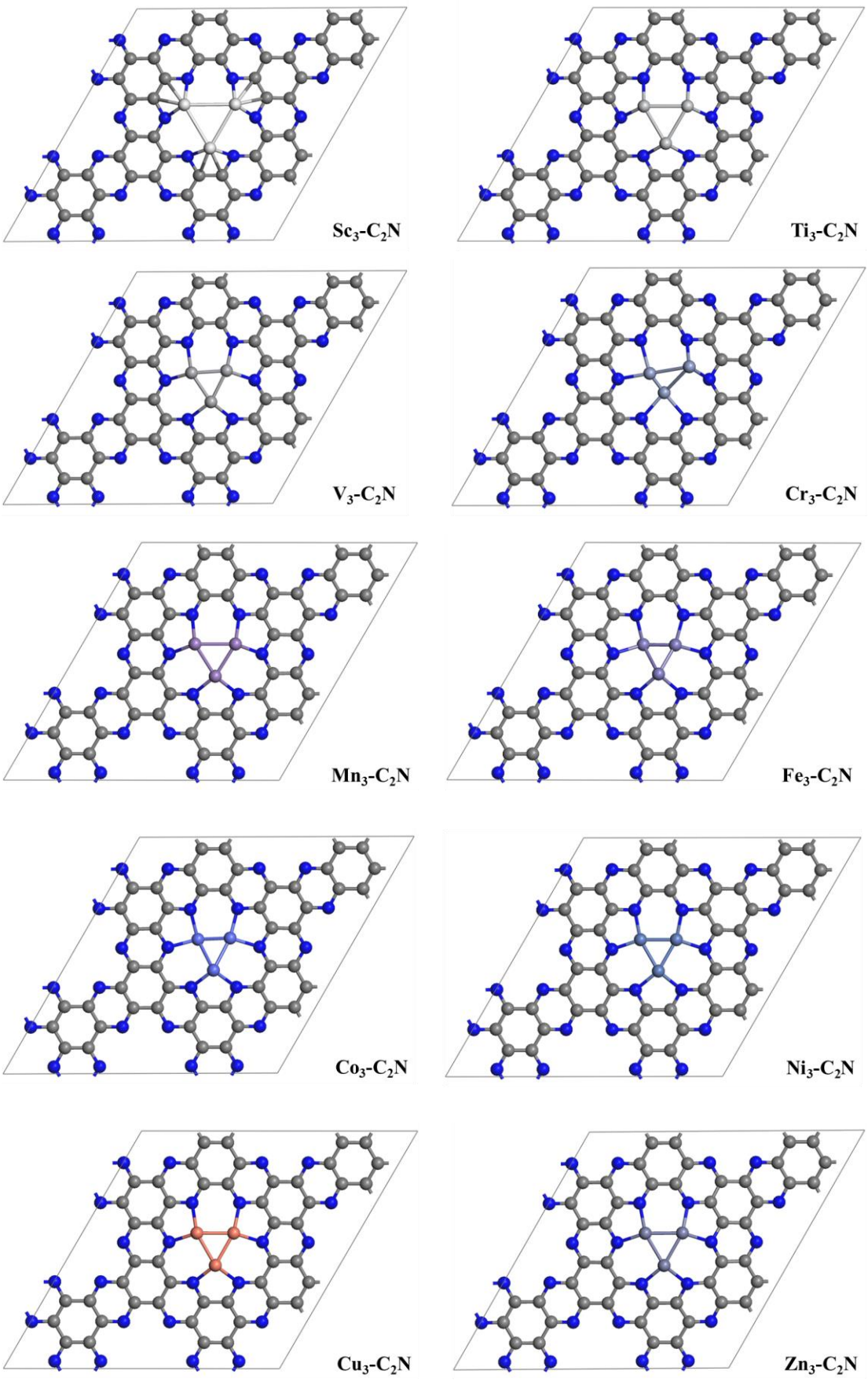


Fig. S1. Optimized structures of M_3-C_2N.