

Supporting information

Fig.S1. (a)-(c) Top and side views of the final Fe-GN4 structures from the molecular dynamics simulation at 700 K, 1000 K and 1300 K. Black, blue and green balls represent the C, N and Fe atom, respectively.

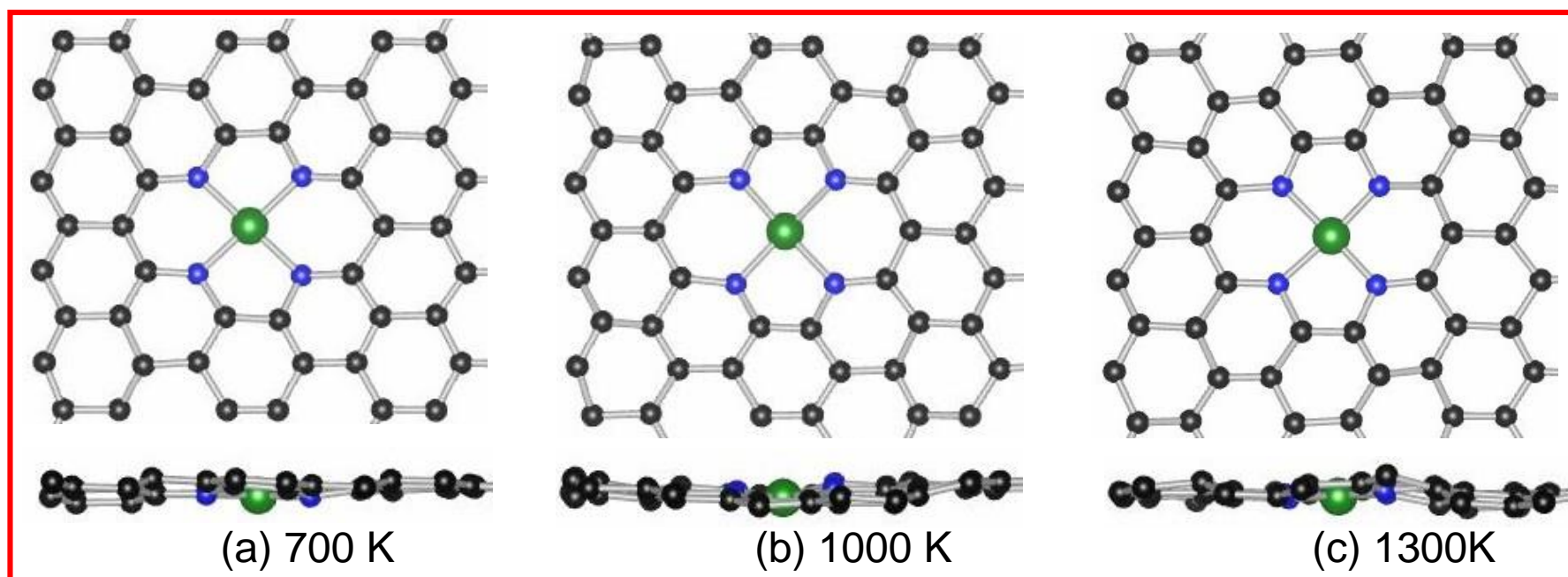


Fig. S1