Cluster Structures and Electronic Properties of FMBeₙ

(FM=Fe, Co, Ni; n=1-12) Clusters

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Fig. S1 Partial density of states (PDOS) of \( s, p, d \) orbits for FMBe\(_n\) \((n=5, 10)\) and Be\(_{n+1}\) \((n=5, 10)\) clusters