supplementary material

Tunable Magnetoelectric Coupling and Electrical Features in an Ultrathin Cr₂Si₂Te₆ / In₂Se₃ Heterostructure

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(a)
(b)
(c)
(c)
(c)
(d)
(d)
(f)

There are six different stacking structures calculated through the DFT calculation, among them, four stacking structures can exist after structural optimization. The energy difference of different stacking structures has been listed in the Table S1, apparently, Stacking I is the most stable structure.

Structure	Ι	II	III	IV
Energy difference (meV)	0	25.7	17.8	17.3

Table. S1. Energy difference of four stacking structure of ultrathin $Cr_2Si_2Te_6$ / In_2Se_3 heterostructure.



Fig. S2. Magneto-crystalline anisotropy energy of 1L in different Hubbard parameter.



Fig. S3. Magneto-crystalline anisotropy energy of 1L Cr₂Si₂Te₆ in different Electric field.