

Supporting Information

Pressure-Induced Conformer Modifications and Electronic Structural Changes in 1, 3, 5-triamino-2, 4, 6-trinitrobenzene (TATB) Upto 20 GPa

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Figure S1. Energy-band diagram of TATB crystal structure at ambient condition

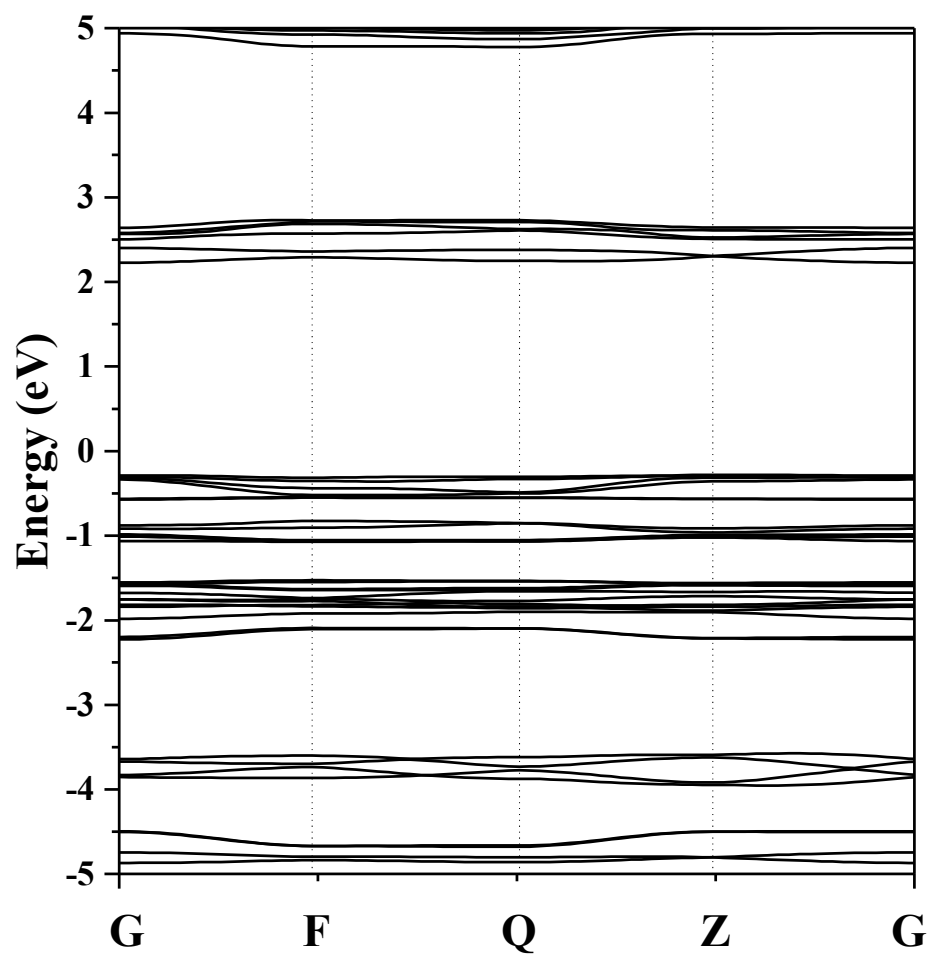


Figure S2. The comparison between experimental and theoretical lattice parameters under compression in this work

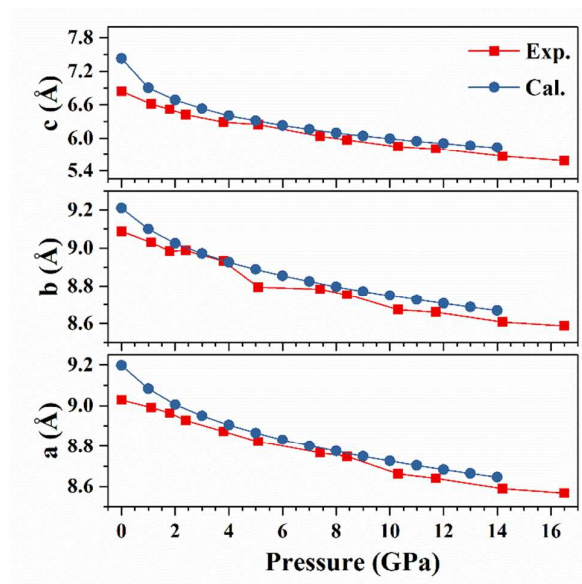


Figure S3. The comparison between theoretical and experimental equation of state in this work

