SKRIPSI

SIMULASI DINAMIKA MOLEKUL STRUKTUR ANATASE-TITANIA (101) TERSULFATASI [TiO_2/SO_4^{2-}] DENGAN METODE *DENSITY-FUNCTIONAL TIGHT-BINDING* (DFTB)

MOLECULAR DYNAMICS SIMULATIONS OF STRUCTURE OF SULFATED ANATASE-TITANIA (101) [TiO_2/SO_4^{2-}] BY DENSITY-FUNCTIONAL TIGHT- BINDING (DFTB)

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