



Formula: C₁₂H₁₄ClNO

MW: 223.7

MDL: MFCD02940828

TNP: TNP00168



LogP: 1.98

LogS: -3.72

Acceptors: 1

Donors: 1

Rotation Bonds: 1

Chiral Centers: 2

N+O: 2

LIPINSKY: 4

IUPAC: 1,2,3,9,9a-pentahydrocyclopenta[2,1-b]quinolin-3-ol, chloride

Smiles: c12c(cccc2)CC2C(=N1)C(CC2)O.Cl