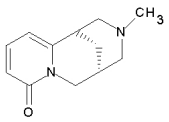

Formula: C12H16N2O

MW: 204.27

MDL: MFCD01459197

TNP:

LogP: 15.12

LogS: -9.84

Acceptors: 1

Donors: 0

Rotation Bonds: 0

Chiral Centers: 2

N+O: 3

LIPINSKY: 3

IUPAC: (1S,9S)-11-methyl-7,11-diazatricyclo[7.3.1.0<2,7>]trideca-2,4-dien-6-one

Smiles: CN1C[C@H]2Cn3c(C(C1)C2)cccc3=O