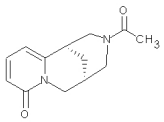

Formula: C13H16N2O2

MW: 232.28

MDL: MFCD01543687

TNP: TNP00329

LogP: 1.57

LogS: -3.21

Acceptors: 2

Donors: 0

Rotation Bonds: 0

Chiral Centers: 2

N+O: 4

LIPINSKY: 4

IUPAC: (9S,1R)-11-acetyl-6-oxo-7,11-diazatricyclo[7.3.1.0<2,7>]trideca-2,4-diene

Smiles: c1(cccc2n1C[C@H]1C[C@@H]2CN(C1)C(=O)C)=O