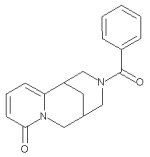

Formula: C18H18N2O2

MW: 294.35

MDL: MFCD02063165

TNP:

LogP: 2.07

LogS: -4.67

Acceptors: 2

Donors: 0

Rotation Bonds: 0

Chiral Centers: 2

N+O: 4

LIPINSKY: 4

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

Smiles: c1ccc(C(N2CC3Cn4c(C(C2)C3)cccc4=O)=O)cc1