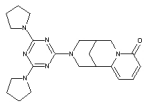

Formula: C22H29N7O

MW: 407.52

MDL: MFCD02579006

TNP:

LogP: 12.62

LogS: -8.71

Acceptors: 1

Donors: 0

Rotation Bonds: 0

Chiral Centers: 2

N+O: 8

LIPINSKY: 3

IUPAC: 11-(4,6-dipyrrolidinyl-1,3,5-triazin-2-yl)-7,11-diazatricyclo[7.3.1.0<2,7>]tri deca-2,4-dien-6-one

Smiles: c1(nc(N2CCCC2)nc(n1)N1CCCC1)N1CC2c3cccc(n3CC(C1)C2)=O