



Formula: C<sub>18</sub>H<sub>20</sub>N<sub>2</sub>O<sub>3</sub>S

MW: 344.43

MDL: MFCD01539409

TNP:



LogP: 5.9

LogS: -5.08

Acceptors: 3

Donors: 0

Rotation Bonds: 0

Chiral Centers: 2

N+O: 5

LIPINSKY: 4

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

Smiles: c12C3CN(CC(C3)Cn1c(ccc2)=O)S(c1ccc(C)cc1)(=O)=O