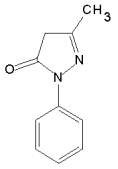

Formula: C10H10N2O

MW: 174.2

MDL NUMBER: MFCD00003138

IUPAC: 3-methyl-1-phenyl-2-pyrazolin-5-one

Smiles: c1ccc(N2C(CC(=N2)C)=O)cc1

ACCEPTORS: 1

DONORS: 0

ROTATION BONDS: 0

N+O: 3

Chiral Centers: 0

LogP: 1.56

LogS: -3.26

LIPINSKI: 4