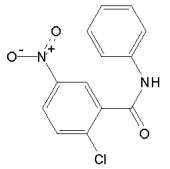

Formula: C13H9ClN2O3

MW: 276.68

MDL NUMBER: MFCD01215270

IUPAC: (2-chloro-5-nitrophenyl)-N-benzamide

Smiles: c1cc(cc(c1Cl)C(Nc1ccccc1)=O)[N+]([O-])=O

Selective PPARgamma antagonist

ACCEPTORS: 3

DONORS: 1

ROTATION BONDS: 1

N+O: 5

Chiral Centers: 0

LogP: 3.48

LogS: -4.06

LIPINSKI: 4