



Formula: C₂₀H₂₁ClO₄

MW: 360.84

CAS: 49562-28-9

MDL NUMBER: MFCD00133314

Smiles: C(C(OC1ccc(C(c2ccc(cc2)Cl)=O)cc1)(C)C)(OC(C)C)=O

THERAPEUTIC CATEGORY: Antihyperlipoproteinemic

ACCEPTORS: 4

DONORS: 0

ROTATION BONDS: 4

N+O: 4

Chiral Centers: 0

LogP: 5.62

LogS: -5.32

LIPINSKI: 4

Synonyms:

2-(4-(4-chlorobenzoyl)phenoxy)-2-methyl-propanoic acid 1-methylethylester; isopropyl(4