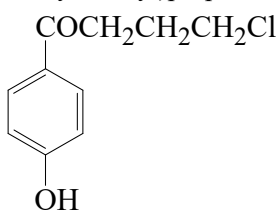




>>Entry Name: 4-Chloro-1-(4-hydroxyphenyl)-1-butanone, 9Cl

Synonym(s): 4-Chloro-4'-hydroxybutyrophenone, 8Cl. 1-Chloro-3-(4-hydroxybenzoyl)propane



CAS Registry Number: 7150-55-2

Molecular Formula: C₁₀H₁₁ClO₂

Molecular Weight: 198.648

Accurate Mass: 198.044757

Percentage Composition: C 60.46%; H 5.58%; Cl 17.85%; O 16.11%

Physical Description: Cryst. (petrol)

Melting Point: Mp 115-117°

Aldrich: C4490-7

>>Derivative: Ac

Molecular Formula: C₁₂H₁₃ClO₃

Molecular Weight: 240.686

Accurate Mass: 240.055322

Percentage Composition: C 59.88%; H 5.44%; Cl 14.73%; O 19.94%

Melting Point: Mp 30-31.5°

>>Derivative: Me ether

Synonym(s): 4-Chloro-1-(4-methoxyphenyl)-1-butanone. 4-Chloro-4'-methoxybutyrophenone

CAS Registry Number: 40877-19-8

Molecular Formula: C₁₁H₁₃ClO₂

Molecular Weight: 212.675

Accurate Mass: 212.060407

Percentage Composition: C 62.12%; H 6.16%; Cl 16.67%; O 15.05%

Physical Description: Plates (EtOAc/petrol)

Melting Point: Mp 31°

Boiling Point: Bp₂ 159-163°

Rare Chemicals Library: S34430-3

References:

van de Westeringh, C. *et al.*, *Ind. Chim. Belg.*, 1960, **25**, 1073-1076, (*synth, Me ether, Ac*)

Kalyanam, N. *et al.*, *Synth. Commun.*, 1988, **18**, 2183-2191, (*synth*)

Uchida, M. *et al.*, *Chem. Pharm. Bull.*, 1989, **37**, 958-961, (*Me ether synth, pmr, ir*)

Chang, S.-J. *et al.*, *Org. Process Res. Dev.*, 2001, **5**, 141-143, (*synth, pmr*)