



>>Entry Name: *N*-[(Diphenylphosphinyl)methyl]-*N*-methylaniline

Synonym(s): *N*-[(Diphenylphosphinyl)methyl]-*N*-methylbenzenamine, 9CI

CAS Registry Number: 76527-75-8

Type of Compound Code(s): AC0310

Ph₂P(O)CH₂NMePh

Molecular Formula: C₂₀H₂₀NOP

Molecular Weight: 321.358

Accurate Mass: 321.128251

Percentage Composition: C 74.75%; H 6.27%; N 4.36%; O 4.98%; P 9.64%

Use/Importance: Reagent for converting ketones into homologues enamines

Physical Description: Cryst. (EtOAc/hexane)

Melting Point: Mp 118-119°

Aldrich: 36741-9

References:

Broekhof, L.N.J.M. *et al.*, *Tet. Lett.*, 1980, **21**, 2671, (*synth, use*)

Van den Gen, A. *et al.*, *ACS Symp. Ser.*, 1981, **171**, 47, (*rev, use*)

Broekhof, N.L.J.M. *et al.*, *Rec. Trav. Chim. (J. R. Neth. Chem. Soc.)*, 1984, **103**, 312, (*synth, ir, pmr, cmr*)

Zorgdrager, J. *et al.*, *Rec. Trav. Chim. (J. R. Neth. Chem. Soc.)*, 1989, **108**, 441, (*synth, ms, pmr, cmr, ir*)