



>>Entry Name: 1-Chloro-4-vinylbenzene

Synonym(s): 1-Chloro-4-ethenylbenzene, 9Cl. *p*-Chlorostyrene, 8Cl. (4-Chlorophenyl)ethylene

CAS Registry Number: 1073-67-2

Extra CAS Registry Numbers(s): 1331-28-8

Structure by analogy with 1-Chloro-2-vinylbenzene, [DXB17-Y](#)

Molecular Formula: C₈H₇Cl

Molecular Weight: 138.596

Accurate Mass: 138.023627

Percentage Composition: C 69.33%; H 5.09%; Cl 25.58%

Melting Point: Mp −15.9°

Boiling Point: Bp 192°. Bp₁₀ 66.3°

Solubility: Insol. H₂O; sol. EtOH, Et₂O, Me₂CO, C₆H₆

Refractive Index: n_D²⁰ 1.5660

Other Data: Coml. samples usually contain polym. inhibitor, e.g. 3,5-di-*tert*-butylcatechol. Undergoes radical polym. Q/e values for copolym., Q 1.33, e −0.64. Rate in radical polym., *k*_p 150 L mol^{−1} s^{−1}, *k*_t 7.7 x 10^{−5} L mol^{−1} s^{−1} (30°)

Aldrich: C7120-3

Fluka: 26360

Sigma: C8423

References:

Aldrich Library of 13C and 1H FT NMR Spectra, 1992, **2**, 140B, (*nmr*)

Aldrich Library of FT-IR Spectra, 1st edn., 1985, **1**, 1019B, (*ir*)

Aldrich Library of FT-IR Spectra: Vapor Phase, 1989, **3**, 951D, (*ir*)

Brooks, L.A., *J.A.C.S.*, 1944, **66**, 1295

Walling, C., *J.A.C.S.*, 1947, **69**, 852

Adv. Chem. Ser., 1955, **15**, 171, (*props*)

Kazuo, N. *et al.*, *Bull. Soc. Chim. Fr.*, 1959, **32**, 771, (*polym*)

Imoto, M. *et al.*, *Makromol. Chem.*, 1965, **86**, 217, (*polym, kinetics*)

Greenley, R.Z., *J. Macromol. Sci., Chem., Part A*, 1980, **14**, 427, (*Q/e values*)