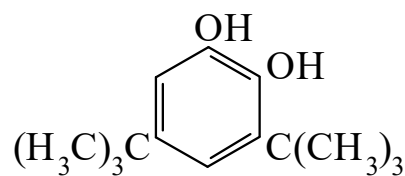




>>Entry Name: 3,5-Di-*tert*-butyl-1,2-benzenediol

Synonym(s): 3,5-Bis(1,1-dimethylethyl)-1,2-benzenediol, 9CI. 3,5-Di-*tert*-butylcatechol



CAS Registry Number: 1020-31-1

Molecular Formula: C₁₄H₂₂O₂

Molecular Weight: 222.327

Accurate Mass: 222.16198

Percentage Composition: C 75.63%; H 9.97%; O 14.39%

Use/Importance: Used as 0.017 M soln. in toluene for extraction-photometric detn. of B ($\lambda_{\max}$ 610 nm, ϵ 105000)

Physical Description: Cryst.

Melting Point: Mp 96-99°

Solubility: Sol. C₆H₆, toluene, CHCl₃

Aldrich: D4580-0

Fluka: 34635

Rare Chemicals Library: S26577-2

References:

Aldrich Library of 13C and 1H FT NMR Spectra, 1992, **2**, 298A, (*nmr*)

Aldrich Library of FT-IR Spectra, 1st edn., 1985, **1**, 1103B, (*ir*)

Aldrich Library of FT-IR Spectra: Vapor Phase, 1989, **3**, 1038C, (*ir*)

Schulze, H. *et al.*, *Annalen*, 1952, **575**, 231, (*synth*)

Oshima, M. *et al.*, *Anal. Chem.*, 1984, **56**, 984, (*detn, B*)