



>>Entry Name: **1,1'-Iminobis-2-propanol**, 9CI

Synonym(s): Diisopropanolamine. DIPA

CAS Registry Number: 110-97-4

H3CCH(OH)CH2NHCH2CH(OH)CH3

Molecular Formula: C₆H₁₅NO₂

Molecular Weight: 133.19

Accurate Mass: 133.110279

Percentage Composition: C 54.11%; H 11.35%; N 10.52%; O 24.02%

Physical Description: Cryst.

Melting Point: Mp 44.5-45.5°

Boiling Point: Bp₇₄₅ 249-250°

Aldrich: 13301-9

Fluka: 14960

>>Derivative: Hydrochloride

CAS Registry Number: 58901-13-6

Physical Description: Cryst. (2-propanol/AcOH)

Melting Point: Mp 79-82°

>>Derivative: Picrate

Physical Description: Prisms (EtOH)

Melting Point: Mp 130-136°

>>Derivative: Bis(4-nitrobenzoyl)

CAS Registry Number: 19207-78-4

Molecular Formula: C₂₀H₂₁N₃O₈

Molecular Weight: 431.401

Percentage Composition: C 55.68%; H 4.91%; N 9.74%; O 29.67%

Physical Description: Cryst. (2-propanol) (as hydrochloride)

Melting Point: Mp 190-192 dec.°

Other Data: CAS no. refers to hydrochloride

>>Derivative: *N*-Nitroso

Synonym(s): 2,2'-Dihydroxy-di-*n*-propylnitrosamine. *N*-Nitrosobis(2-hydroxypropyl)amine. Diisopropanolnitrosamine

CAS Registry Number: 53609-64-6

Molecular Formula: C₆H₁₄N₂O₃

Molecular Weight: 162.188

Accurate Mass: 162.100443

Percentage Composition: C 44.43%; H 8.70%; N 17.27%; O 29.59%

Boiling Point: Bp_{0.01} 95°

Refractive Index: n_D²¹ 1.4826

Hazard & Toxicity: Carcinogen

References:

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Kretov, A.E. *et al.*, *Zh. Obshch. Khim.*, 1962, **32**, 471; *CA*, **58**, 1337d, (*synth*)

Krueger, F.W. *et al.*, *Naturwissenschaften*, 1974, **61**, 328, (*N-nitroso, synth, tox*)

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