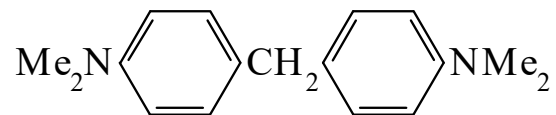




>>Entry Name: Bis(4-dimethylaminophenyl)methane

Synonym(s): 4,4'-Methylenebis[*N,N*-dimethylbenzenamine], 9Cl. 4,4'-Methylenebis[*N,N*-dimethylaniline], 8Cl. 4,4'-Tetramethyldiaminodiphenylmethane. Methane base. Michler's base. Tetra base. Arnold's base



CAS Registry Number: 101-61-1

Molecular Formula: C₁₇H₂₂N₂

Molecular Weight: 254.374

Accurate Mass: 254.178298

Percentage Composition: C 80.27%; H 8.72%; N 11.01%

Use/Importance: Used as 2% soln. in Me₂CO to give colour reaction with [Fe(CN)₆]³⁽⁻⁾

Physical Description: Leaflets or plates (EtOH or petrol)

Melting Point: Mp 91°

Boiling Point: Bp 390°. Bp₃ 182-185°

Solubility: Sol. Me₂CO

Hazard & Toxicity: Exp. carcinogen. LD₅₀ (mus, orl) 3160 mg/kg

Aldrich: M4445-1

Fluka: 87802

Rare Chemicals Library: S43663-1

Sigma: M6144

>>Derivative: Dipicrate

Melting Point: Mp 178°

>>Derivative: Methiodide (1:2)

Physical Description: Yellow leaflets

Melting Point: Mp 214 dec.°

Other Data: Becomes green at 193°

References:

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Mangini, A. *et al.*, *J.C.S.*, 1956, 4954, (*uv*)

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Gaspariř, J. *et al.*, *Coll. Czech. Chem. Comm.*, 1959, **24**, 1943, (*chromatog*)

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