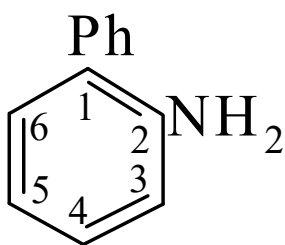




>>Entry Name: 2-Aminobiphenyl

Synonym(s): [1,1'-Biphenyl]-2-amine, 9CI



CAS Registry Number: 90-41-5

Molecular Formula: C₁₂H₁₁N

Molecular Weight: 169.226

Accurate Mass: 169.089149

Percentage Composition: C 85.17%; H 6.55%; N 8.28%

Use/Importance: Fluorogenic reagent for the detn. of carbonyl compds.

Physical Description: Cryst. (EtOH aq.)

Melting Point: Mp 49-50°

Boiling Point: Bp 299°. Bp₁₅ 170°. Bp₅ 145-148°

Other Data: Steam-volatile

Hazard & Toxicity: Autoignition temp. 450°. LD₅₀ (rat, orl) 2340 mg/kg

Aldrich: A4240-9

Fluka: 7110

Sigma: A1129

>>Derivative: Hydrochloride

CAS Registry Number: 2185-92-4

Physical Description: Cryst.

Melting Point: Mp 180°

Hazard & Toxicity: Exp. carcinogen

Rare Chemicals Library: S41293-7

>>Derivative: *N*-Formyl

Synonym(s): *N*-[1,1'-Biphenyl]-2-ylformamide, 9CI

CAS Registry Number: 5346-21-4

Molecular Formula: C₁₃H₁₁NO

Molecular Weight: 197.236

Accurate Mass: 197.084064

Percentage Composition: C 79.17%; H 5.62%; N 7.10%; O 8.11%

Physical Description: Needles (EtOH aq.)

Melting Point: Mp 75°

Rare Chemicals Library: S84102-1

>>Derivative: *N*-Ac

Synonym(s): 2-Acetamidobiphenyl

CAS Registry Number: 2113-47-5

Molecular Formula: C₁₄H₁₃NO

Molecular Weight: 211.263

Accurate Mass: 211.099714

Percentage Composition: C 79.59%; H 6.20%; N 6.63%; O 7.57%

Physical Description: Needles (EtOH aq. or petrol)

Melting Point: Mp 121°

Boiling Point: Bp 355°

Rare Chemicals Library: S81126-2