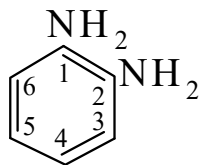




>>Entry Name: **1,2-Benzenediamine**, 9CI  
Synonym(s): *o*-Phenylenediamine, 8CI. 1,2-Diaminobenzene



CAS Registry Number: 95-54-5  
Related CAS Registry Numbers(s): 74710-09-1  
Extra CAS Registry Numbers(s): 25265-76-3

Molecular Formula: C<sub>6</sub>H<sub>8</sub>N<sub>2</sub>  
Molecular Weight: 108.143  
Accurate Mass: 108.068748  
Percentage Composition: C 66.64%; H 7.46%; N 25.90%  
Source/Synthesis: Manuf. by Fe-HCl redn. of 1,2-dinitrobenzene or 1-nitroaniline  
Use/Importance: Used as 0.2% aq. soln. for photometric detn. of Br<sup>(-)</sup>, Se, Pt, V. Intermed. for heterocyclic synth., esp. of quinoxalines. Used in manuf. of antioxidants, photographic chemicals and polyamide fibres. Chromogenic substrate used in conjunction with horseradish peroxidase  
Physical Description: Leaflets (H<sub>2</sub>O)  
Melting Point: Mp 104°  
Boiling Point: Bp 256-258°  
Solubility: Sol. hot H<sub>2</sub>O, EtOH, Et<sub>2</sub>O, CHCl<sub>3</sub>  
pKa Value: pK<sub>a1</sub> 4.65; pK<sub>a2</sub> 1.86 (20°, H<sub>2</sub>O)  
Other Data: Electrochem. homopolym. gives an electroactive ladder polymer containing phenazine rings  
**Hazard & Toxicity:** Eye and skin irritant and may cause dermatitis. LD<sub>50</sub> (rat, orl) 1070 mg/kg

Aldrich: 27578-6  
Fluka: 78412  
Sigma: O6882

>>Derivative: Hydrochloride (1:2)

CAS Registry Number: 615-28-1  
Physical Description: Pink-white cryst. (MeOH)  
Melting Point: Mp 290 dec.°

Aldrich: 23823-6  
Fluka: 78440  
Rare Chemicals Library: S57083-4  
Sigma: P8787

>>Derivative: Picrate

Physical Description: Orange monoclinic cryst.  
Melting Point: Mp 214° (dec.)

>>Derivative: *N,N'*-Diformyl  
Synonym(s): *N,N'*-1,2-Phenylenebisformamide, 9CI. 1,2-Bis(formamido)benzene

CAS Registry Number: 31354-51-5  
Molecular Formula: C<sub>8</sub>H<sub>8</sub>N<sub>2</sub>O<sub>2</sub>  
Molecular Weight: 164.163  
Accurate Mass: 164.058578  
Percentage Composition: C 58.53%; H 4.91%; N 17.06%; O 19.49%  
Melting Point: Mp 155-157 dec.°

>>Derivative: *N*-Ac  
Synonym(s): *N*-(2-Aminophenyl)acetamide, 9CI. 2-Aminoacetanilide

CAS Registry Number: 34801-09-7  
Molecular Formula: C<sub>8</sub>H<sub>10</sub>N<sub>2</sub>O  
Molecular Weight: 150.18  
Accurate Mass: 150.079313  
Percentage Composition: C 63.98%; H 6.71%; N 18.65%; O 10.65%  
Physical Description: Plates (C<sub>6</sub>H<sub>6</sub>)  
Melting Point: Mp 132° (145°)