



>>Entry Name: 3-Fluorophenylalanine, 9CI
Synonym(s): 2-Amino-3-(3-fluorophenyl)propanoic acid

CAS Registry Number: 456-88-2
Structure by analogy with 2-Fluorophenylalanine, [CJB83-K](#)

Molecular Formula: C₉H₁₀FNO₂
Molecular Weight: 183.182
Accurate Mass: 183.069557
Percentage Composition: C 59.01%; H 5.50%; F 10.37%; N 7.65%; O 17.47%
Hazard & Toxicity: LD₅₀ (mus, ipr) 5.9 mg/kg

>>Variant: (R)-form
Synonym(s): D-form

Molecular Formula: C₉H₁₀FNO₂
Molecular Weight: 183.182
Accurate Mass: 183.069557
Percentage Composition: C 59.01%; H 5.50%; F 10.37%; N 7.65%; O 17.47%
Physical Description: Cryst. (EtOH aq.)
Melting Point: Mp 230-234°
Optical Rotation: $[\alpha]_D^{25} +22$ (c, 2 in H₂O)

>>Derivative: N-Benzyloxycarbonyl

Melting Point: Mp 126-127°
Optical Rotation: $[\alpha]_D^{18} +7.35$ (c, 1 in MeOH)

>>Variant: (S)-form
Synonym(s): L-form

CAS Registry Number: 19883-77-3
Molecular Formula: C₉H₁₀FNO₂
Molecular Weight: 183.182
Accurate Mass: 183.069557
Percentage Composition: C 59.01%; H 5.50%; F 10.37%; N 7.65%; O 17.47%
Physical Description: Cryst. (EtOH aq.)
Melting Point: Mp 239-243° (235-239°)
Optical Rotation: $[\alpha]_D^{18} -25.2$ (c, 1 in H₂O)
Fluka: 47306

>>Derivative: N-Benzyloxycarbonyl

Physical Description: Cryst. (diisopropyl ether)
Melting Point: Mp 124°
Optical Rotation: $[\alpha]_D^{20} -6.2$ (c, 1 in MeOH)

>>Variant: (±)-form

CAS Registry Number: 2629-54-1
Molecular Formula: C₉H₁₀FNO₂
Molecular Weight: 183.182
Accurate Mass: 183.069557
Percentage Composition: C 59.01%; H 5.50%; F 10.37%; N 7.65%; O 17.47%
Melting Point: Mp 240-250°. Mp 262-263°
Aldrich: 21943-6
Fluka: 47310
Sigma: F5126

>>Derivative: Hydrochloride

Melting Point: Mp 248-251 dec.°