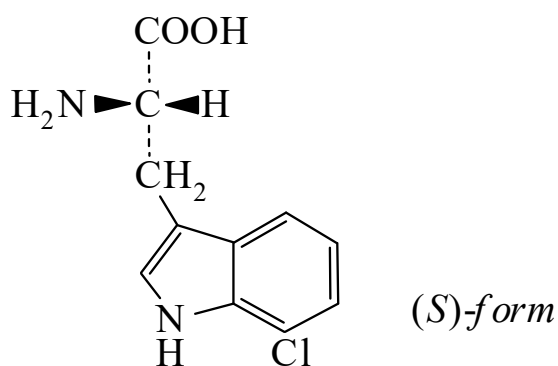




>>Entry Name: 7-Chlorotryptophan, 9Cl

Synonym(s): 2-Amino-3-(7-chloro-1*H*-indol-3-yl)propanoic acid



Extra CAS Registry Numbers(s): 153-97-9

Molecular Formula: C₁₁H₁₁ClN₂O₂

Molecular Weight: 238.673

Accurate Mass: 238.050905

Percentage Composition: C 55.36%; H 4.65%; Cl 14.85%; N 11.74%; O 13.41%

Biological Source: Either enantiomer may serve as a precursor of Pyrrolnitrin [CKL72-L](#)

>>Variant: (*R*)-form

Synonym(s): D-form

CAS Registry Number: 75102-74-8

Molecular Formula: C₁₁H₁₁ClN₂O₂

Molecular Weight: 238.673

Accurate Mass: 238.050905

Percentage Composition: C 55.36%; H 4.65%; Cl 14.85%; N 11.74%; O 13.41%

Melting Point: Mp 275 dec.°

Optical Rotation: $[\alpha]_{\text{D}}^{23} +32.5$ (c, 0.04 in H₂O)

>>Variant: (*S*)-form

Synonym(s): L-form

CAS Registry Number: 73945-46-7

Molecular Formula: C₁₁H₁₁ClN₂O₂

Molecular Weight: 238.673

Accurate Mass: 238.050905

Percentage Composition: C 55.36%; H 4.65%; Cl 14.85%; N 11.74%; O 13.41%

Melting Point: Mp 270 dec.°

Optical Rotation: $[\alpha]_{\text{D}}^{23} -29.7$ (c, 0.06 in H₂O)

References:

van Pée, K.-H. *et al.*, *Angew. Chem., Int. Ed.*, 1980, **19**, 828, (*biosynth*)

van Pée, K.-H. *et al.*, *Annalen*, 1981, 233, (*synth*)

Lee, M. *et al.*, *J. Het. Chem.*, 1994, **31**, 711, (*pmr, cmr*)