



>>Entry Name: Benzyloxycarbonylcystine

Synonym(s): *N,N'*-Bis[(Phenylmethoxy)carbonyl]cystine, 9Cl. *Z*-Cystine. (*Z*-CysOH)₂

PhCH₂OOCNHCH(COOH)CH₂SSCH₂CH(COOH)NHCOOCH₂Ph

Molecular Formula: C₂₂H₂₄N₂O₈S₂

Molecular Weight: 508.572

Accurate Mass: 508.097408

Percentage Composition: C 51.96%; H 4.76%; N 5.51%; O 25.17%; S 12.61%

>>Variant: L-form

CAS Registry Number: 6968-11-2

Molecular Formula: C₂₂H₂₄N₂O₈S₂

Molecular Weight: 508.572

Accurate Mass: 508.097408

Percentage Composition: C 51.96%; H 4.76%; N 5.51%; O 25.17%; S 12.61%

Melting Point: Mp 119-120°

Optical Rotation: [α]_D −51.6 (c, 1 in DMF)

Fluka: 96080

>>Derivative: Dicyclohexylamine salt

Synonym(s): (*Z*-CysOH.Dcha)₂

CAS Registry Number: 38595-81-2

Melting Point: Mp 190-191°

Optical Rotation: [α]_D −51.8 (c, 1 in DMF)

>>Derivative: Dihydrazide

Synonym(s): (*Z*-CysNHNH₂)₂

Molecular Formula: C₂₂H₂₈N₆O₆S₂

Molecular Weight: 536.632

Accurate Mass: 536.151174

Percentage Composition: C 49.24%; H 5.26%; N 15.66%; O 17.89%; S 11.95%

Melting Point: Mp 175-176°

>>Derivative: Bispentachlorophenyl ester

Synonym(s): (*Z*-CysOPcp)₂

CAS Registry Number: 13673-50-2

Melting Point: Mp 215-217°

Optical Rotation: [α]_D −55.6 (c, 0.63 in CHCl₃)

>>Derivative: Disuccinimido ester

Synonym(s): (*Z*-CysOSu)₂

Melting Point: Mp 112-113°

References:

King, L.C. *et al.*, *J.A.C.S.*, 1952, **74**, 5499, (*synth*)

Zervas, L. *et al.*, *J.A.C.S.*, 1959, **81**, 1729, (*hydrazide*)

Klieger, E. *et al.*, *Annalen*, 1961, **640**, 157, (*synth*)

Kovacs, J. *et al.*, *J.O.C.*, 1967, **32**, 3696, (*pentachlorophenyl*)

Gorlenko, V.A. *et al.*, *J. Gen. Chem. USSR (Engl. Transl.)*, 1970, **40**, 1623, (*succinimido*)