



>>Entry Name: Benzyloxycarbonylcysteine

Synonym(s): *N*-[(Phenylmethoxy)carbonyl]cysteine. Z-Cysteine. Z-CysOH

PhCH₂OOCNHCH(COOH)CH₂SH

Molecular Formula: C₁₁H₁₃NO₄S

Molecular Weight: 255.294

Accurate Mass: 255.056529

Percentage Composition: C 51.75%; H 5.13%; N 5.49%; O 25.07%; S 12.56%

>>Variant: L-form

CAS Registry Number: 53907-29-2

Molecular Formula: C₁₁H₁₃NO₄S

Molecular Weight: 255.294

Accurate Mass: 255.056529

Percentage Composition: C 51.75%; H 5.13%; N 5.49%; O 25.07%; S 12.56%

Melting Point: Mp 163-165°

Optical Rotation: [α]_D −134 (DMF)

>>Derivative: *S*-Acetamidomethyl

Synonym(s): Z-Cys(Acm)OH

CAS Registry Number: 54784-68-8

Physical Description: Oil

>>Derivative: *S*-Acetamidomethyl, succinimido ester

Synonym(s): Z-Cys(Acm)OSu

CAS Registry Number: 38004-20-5

Melting Point: Mp 152-154°

>>Derivative: *S*-Benzyl, dicyclohexylamine salt

Synonym(s): Z-Cys(Bzl)OH.Dcha

CAS Registry Number: 3257-18-9

Melting Point: Mp 138-139°

Optical Rotation: [α]_D −3.9 (c, 2.01 in EtOH)

Aldrich: 12101-0

Sigma: C5876

>>Derivative: *S*-Benzyl, hydrazide

Synonym(s): Z-Cys(Bzl)NHNH₂

CAS Registry Number: 24164-46-3

Molecular Formula: C₁₈H₂₁N₃O₃S

Molecular Weight: 359.448

Accurate Mass: 359.130362

Percentage Composition: C 60.15%; H 5.89%; N 11.69%; O 13.35%; S 8.92%

Melting Point: Mp 133-135°

Optical Rotation: [α]_D −14.3 (c, 1 in EtOH)

>>Derivative: *S*-Benzyl, 4-nitrophenyl ester

Synonym(s): Z-Cys(Bzl)ONp

CAS Registry Number: 3401-37-4

Melting Point: Mp 93-94°

Optical Rotation: [α]_D −43 (c, 2 in DMF)

Sigma: C6001

>>Derivative: *S*-Benzyl, succinimido ester

Synonym(s): Z-Cys(Bzl)OSu

CAS Registry Number: 3401-57-8

Melting Point: Mp 90-91°

Optical Rotation: [α]_D −58.5 (c, 1.8 in dioxan)