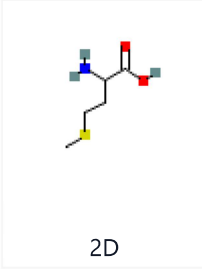



COMPOUND SUMMARY

Racemethionine

PubChem CID	876
Structure	 2D  3D
Chemical Safety	Laboratory Chemical Safety Summary (LCSS) Datasheet
Molecular Formula	C₅H₁₁NO₂S CH ₃ S(CH ₂) ₂ CH(NH ₂)COOH
Synonyms	DL-METHIONINE 59-51-8 Racemethionine methionine Acimetion View More...
Molecular Weight	149.21 g/mol <i>Computed by PubChem 2.2 (PubChem release 2021.10.14)</i>
Dates	Create: 2004-09-16 Modify: 2024-03-30
Description	Methionine is a sulfur-containing amino acid that is butyric acid bearing an amino substituent at position 2 and a methylthio substituent at position 4. It has a role as an Escherichia coli metabolite, a Saccharomyces cerevisiae metabolite, a plant metabolite, a Daphnia magna metabolite and an algal metabolite. It is an alpha-amino acid and a sulfur-containing amino acid. It is functionally related to a butyric acid . It is a conjugate base of a

methioninium. It is a conjugate acid of a **methioninate**. It is a tautomer of a **methionine** zwitterion.

▶ [ChEBI](#)

A preparation of **methionine** that includes a mixture of **D-methionine** and **L-methionine** isomers.

▶ [DrugBank](#)

DL-Methionine is a natural product found in **Drosophila melanogaster**, **Mycoplasma gallisepticum**, and **other organisms** with data available.

▶ [LOTUS - the natural products occurrence database](#)

See also: ... [View More ...](#)

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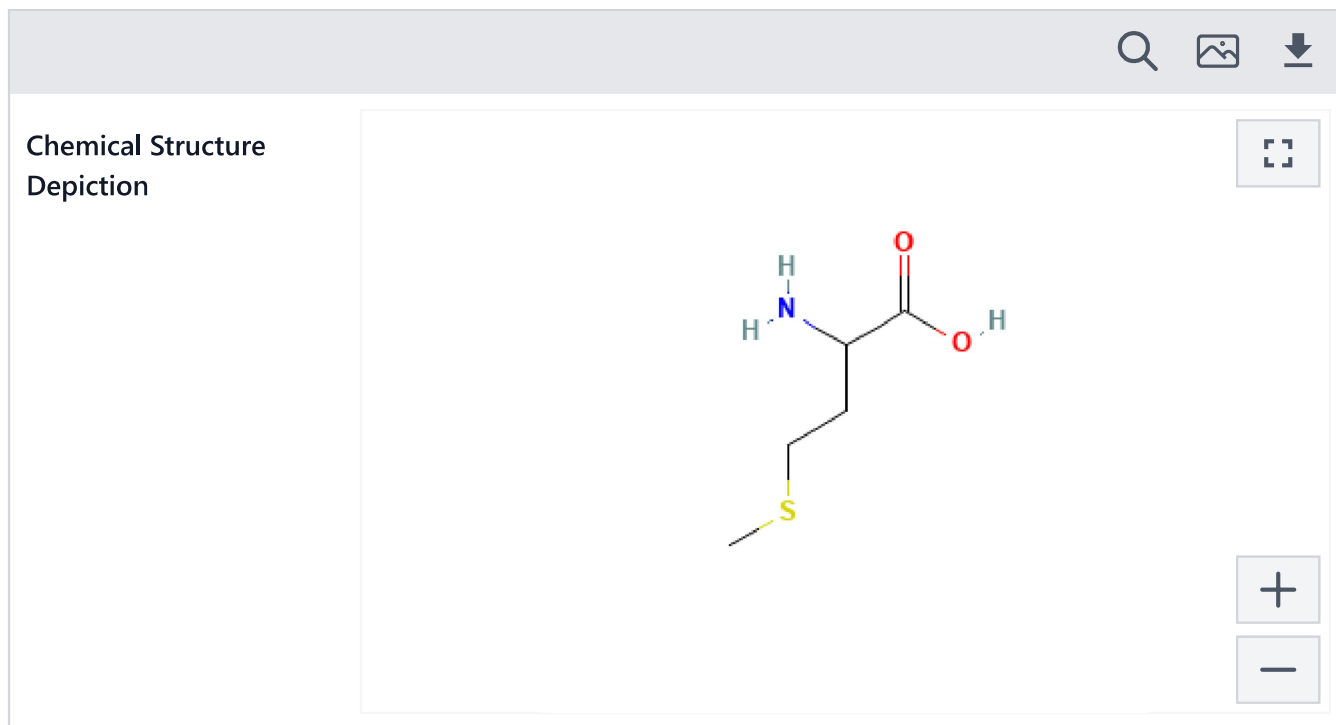


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1 Structures



1.1 2D Structure

[PubChem](#)

1.2 3D Conformer



► [PubChem](#)

2 Biologic Description



SVG Image



IUPAC Condensed	H-DL-Met-OH
Sequence	M
PLN	H-M-OH
HELM	PEPTIDE1{(M,[dM])}\$\$\$\$V2.0
IUPAC	DL-methionine

► [PubChem](#)

3 Names and Identifiers



3.1 Computed Descriptors



3.1.1 IUPAC Name



2-amino-4-methylsulfanylbutanoic acid

Computed by Lexichem TK 2.7.0 (PubChem release 2021.10.14)

▶ [PubChem](#)

3.1.2 InChI



InChI=1S/C5H11NO2S/c1-9-3-2-4(6)5(7)8/h4H,2-3,6H2,1H3,(H,7,8)

Computed by InChI 1.0.6 (PubChem release 2021.10.14)

▶ [PubChem](#)

3.1.3 InChIKey



FFEARJCKVFRZRR-UHFFFAOYSA-N

Computed by InChI 1.0.6 (PubChem release 2021.10.14)

▶ [PubChem](#)

3.1.4 Canonical SMILES



CSCCC(C(=O)O)N

Computed by OEChem 2.3.0 (PubChem release 2021.10.14)

▶ [PubChem](#)

3.2 Molecular Formula



C₅H₁₁NO₂S

CH₃S(CH₂)₂CH(NH₂)COOH