

L-Tryptophan

Other names:	(+)-tryptophan (-)-Tryptophan (R)-.alpha.-amino-3-indolepropionic acid (R)-tryptophan (S)-Tryptophan (S)-«alpha»-Amino-1H-indole-3-propanoic acid 1H-Indole-3-alanine 1H-Indole-3-alanine, (S)- 1H-Indole-3-propanoic acid, «alpha»-amino-, (S)- 2-Amino-3-indolylpropanoic acid 3-Indol-3-ylalanine ALPHA-AMINOINDOLE-3-PROPIONIC ACID Alanine, 3-indol-3-yl- D-tryptophan EH 121 Indole-3-alanine Indole-3-propionic acid, «alpha»-amino- L-(-)-Tryptophan L-.alpha.-amino-3-indolepropionic acid L-2-amino-3-indolepropanoic acid L-Alanine, 3-(1H-indol-3-yl)- L-Trp L-Tryptofan L-Tryptophane L-«alpha»-amino-3-indolepropionic acid L-«alpha»-aminoindole-3-propionic acid L-«beta»-3-indolylalanine NCI-C01729 NSC 13119 Pacitron Propionic acid, 2-amino-3-indol-3-yl- Trp Tryptophan Tryptophan, L- Tryptophane propanoic acid, 2-amino-3-indole-, L- «alpha»'-Amino-3-indolepropionic acid «alpha»-Aminoindole-3-propionic acid
Inchi:	InChI=1S/C11H12N2O2/c12-9(11(14)15)5-7-6-13-10-4-2-1-3-8(7)10/h1-4,6,9,13H,5,12H
InchiKey:	QIVBCDIJIAJPQS-SECBINFHSA-N

Formula: C11H12N2O2
SMILES: NC(Cc1c[nH]c2ccccc12)C(=O)O
Mol. weight [g/mol]: 204.23
CAS: 73-22-3

Physical Properties

Property code	Value	Unit	Source
affp	948.90	kJ/mol	NIST Webbook
affp	927.20	kJ/mol	NIST Webbook
affp	945.60	kJ/mol	NIST Webbook
basg	915.00	kJ/mol	NIST Webbook
chs	-5628.32 ± 0.84	kJ/mol	NIST Webbook
log10ws	-2.36		Crippen Method
logp	0.640		Crippen Method
mcvol	154.330	ml/mol	McGowan Method
ss	251.04	J/mol×K	NIST Webbook

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cps	238.15	J/mol×K	298.15	NIST Webbook
hvapt	184.40	kJ/mol	450.15	Enthalpy of sublimation of natural aromatic amino acids determined by Knudsen's effusion mass spectrometric method
rhos	1080.00	kg/m3	298.15	Thermodynamic properties of L-tryptophan

Sources

Viscosities of L-Histidine/L-Glutamic Acid/L-Tryptophan/Glycylglycine+2 M Aqueous KCl/KNO₃ Solutions at T =(298.15 to 323.15)K: <https://www.doi.org/10.1007/s10765-011-0996-9>

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C73223&Units=SI
Saturated Solubility and Thermodynamic Evaluation of McGowan Method	https://www.doi.org/10.1021/acs.jced.9b00562
Enthalpy of sublimation of natural aromatic amino acids determined by Kihara's thermodynamic properties of extrapolation method:	http://link.springer.com/article/10.1007/BF02311772
Ultrasonic Velocities and Densities of L-Histidine or L-Glutamic Acid or Cysteine Mixture Glycylglycine + 2 mol/L Aqueous KCl or KNO₃ Solutions from 298.15 to 323.15 K:	https://www.cheric.org/research/kdb/hcprop/showprop.php?cmpid=1490
L-Tyrosine in Several Organic Solvents: Experimental solubility, enthalpy and thermodynamic analysis of biologically soluble D,L-tryptophan in aqueous α-Amino Acids Water/Ethanol/N,N-dimethylformamide binary mixtures.	https://www.doi.org/10.1016/j.jct.2014.02.003
Temperature of Maximum Density for Binary Aqueous Solutions of Five Amino acids.	https://www.doi.org/10.1016/j.jct.2016.09.041
PAMAM-NH₂ G4 dendrimer with selected natural amino acids in aqueous solutions:	https://www.doi.org/10.1021/je900199j
	https://www.chemeo.com/doc/models/crippen_log10ws
	https://www.doi.org/10.1021/acs.jced.9b00258
	https://www.doi.org/10.1016/j.jct.2018.08.018
	https://www.doi.org/10.1021/acs.jced.7b00486
	https://www.doi.org/10.1016/j.fluid.2011.09.028
	https://www.doi.org/10.1021/acs.jced.9b00752
	https://www.doi.org/10.1016/j.jct.2013.10.022

Legend

affp:	Proton affinity
basg:	Gas basicity
chs:	Standard solid enthalpy of combustion
cps:	Solid phase heat capacity
hvapt:	Enthalpy of vaporization at a given temperature
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
rhos:	Solid Density
ss:	Solid phase molar entropy at standard conditions

Latest version available from:

<https://www.chemeo.com/cid/32-606-3/L-Tryptophan.pdf>

Generated by Cheméo on 2024-04-25 15:45:18.905875098 +0000 UTC m=+16349167.826452410.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.