

N-(N-Glycylglycyl)glycine

Other names:

Diglycylglycine
Glycine, N-(N-glycylglycyl)-
Glycylglycylglycine
Triglycine

Inchi: InChI=1S/C6H11N3O4/c7-1-4(10)8-2-5(11)9-3-6(12)13/h1-3,7H2,(H,8,10)(H,9,11)(H,12,13)

InchiKey: XKUKSGPZAADMRA-UHFFFAOYSA-N

Formula: C6H11N3O4

SMILES: NCC(O)=NCC(O)=NCC(=O)O

Mol. weight [g/mol]: 189.17

CAS: 556-33-2

Physical Properties

Property code	Value	Unit	Source
hf	-557.79	kJ/mol	Joback Method
hvap	103.16	kJ/mol	Joback Method
log10ws	1.17		Crippen Method
logp	-1.057		Crippen Method
mcvol	135.920	ml/mol	McGowan Method
pc	3916.03	kPa	Joback Method
tb	892.74	K	Joback Method
tc	1098.97	K	Joback Method

Sources

Enthalpic Interaction Coefficients of N-(N-Glycylglycyl)glycine in Aqueous Solutions of Organic Acids and Glycine Salts with Organic Sodium Sulfate and Glycylglycylglycine

<https://www.doi.org/10.1021/je900849b>

Thermodynamic properties of peptide solutions 20. Partial molar volumes and isobaric heat capacities for some tripeptides of sequence gly-X-gly (X = Gly, Ala, Leu, Asp, Glu, and Tyr) in aqueous solutions containing model compounds of amino acids and their pic-salts at 0.1 mol/kg with the osmolyte trimethylamine N-oxide in aqueous solutions

<https://www.doi.org/10.1016/j.jct.2003.09.010>

<https://www.doi.org/10.1016/j.tca.2009.02.017>

<http://pubs.acs.org/doi/abs/10.1021/ci9903071>

https://www.chemeo.com/doc/models/crippen_log10ws

<https://www.doi.org/10.1016/j.jct.2016.03.032>

https://en.wikipedia.org/wiki/Joback_method

<https://www.doi.org/10.1016/j.jct.2006.11.014>

<https://www.doi.org/10.1016/j.jct.2011.12.029>

<https://www.doi.org/10.1021/je0600754>

Studies on the interactions of diglycine and triglycine with polyethylene glycol (PEG) in aqueous solutions by density and ultrasound speed measurements: Thermodynamics of the interactions of some amino acids and peptides with McGowan's Method and tetradecyltrimethylammonium bromide. Partial molar volumes, expansibilities and compressibilities of glyglyglycine interactions with glucose and fructose pentoses with the drug pentoxifyllin in aqueous solution compared with glucose in aqueous solution. A new approach: temperatures of some inorganic amino acids, peptides and protein with aqueous 3-chloro-1,2-propanediol and 3-chloro-1-propanol: Biophysical studies:

<https://www.doi.org/10.1016/j.jct.2012.10.018>
<http://webbook.nist.gov/cgi/cbook.cgi?ID=C556332&Units=SI>
<https://www.doi.org/10.1016/j.jct.2013.11.001>
<http://link.springer.com/article/10.1007/BF02311772>
<https://www.doi.org/10.1016/j.tca.2010.11.013>
<https://www.doi.org/10.1016/j.jct.2012.05.009>
<https://www.doi.org/10.1016/j.tca.2010.05.016>
<https://www.doi.org/10.1016/j.jct.2010.11.015>

Legend

hf:	Enthalpy of formation at standard conditions
hvpv:	Enthalpy of vaporization at standard conditions
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature

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