

allo-L-isoleucine

Other names:	(2S,3S)-.alpha.-amino-.beta.-methyl-n-valeric acid (2S,3S)-.alpha.-amino-.beta.-methylvaleric acid (2S,3S)-2-amino-3-methylpentanoic acid (S)-isoleucine (S,S)-isoleucine .alpha.-amino-.beta.-methylvaleric acid 2-amino-3-methylpentanoic acid 2-amino-3-methylvaleric acid 2S,3S-isoleucine L-(+)-isoleucine L-isoleucine L-norvaline, 3-methyl-, erythro- [S-(R*,R*)]-2-amino-3-methylpentanoic acid erythro-L-isoleucine isoleucine isoleucine, L- pentanoic acid, 2-amino-3-methyl-, [S-(R*,R*)]-
Inchi:	InChI=1S/C6H13NO2/c1-3-4(2)5(7)6(8)9/h4-5H,3,7H2,1-2H3,(H,8,9)
InchiKey:	AGPKZVBTJJNPAG-UHFFFAOYSA-N
Formula:	C6H13NO2
SMILES:	CCC(C)C(N)C(=O)O
Mol. weight [g/mol]:	131.17
CAS:	3107-04-8

Physical Properties

Property code	Value	Unit	Source
gf	-204.53	kJ/mol	Joback Method
hf	-408.75	kJ/mol	Joback Method
hfus	15.13	kJ/mol	Joback Method
hvap	62.24	kJ/mol	Joback Method
log10ws	-0.74		Crippen Method
logp	0.444		Crippen Method
mcvol	112.820	ml/mol	McGowan Method
pc	4077.71	kPa	Joback Method
tb	554.38	K	Joback Method
tc	743.19	K	Joback Method
tf	321.39	K	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	275.71	J/molxK	554.38	Joback Method
cpg	285.39	J/molxK	585.85	Joback Method
cpg	294.59	J/molxK	617.32	Joback Method
cpg	303.34	J/molxK	648.79	Joback Method
cpg	311.64	J/molxK	680.26	Joback Method
cpg	319.50	J/molxK	711.72	Joback Method
cpg	326.94	J/molxK	743.19	Joback Method
hsubt	120.10 ± 0.80	kJ/mol	455.00	NIST Webbook

Sources

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<http://pubs.acs.org/doi/abs/10.1021/ci9903071>
- Effect of concentration and temperature on apparent molar expansivities of some of the alpha-amino hydrochlorides (Glycine/L-alanine) in aqueous solutions. Solvation in the water-urea-DMSO and DMSO-ethanol binary systems. Effect of the hydrophobicity of the amino acid on the apparent molar expansivities in aqueous solutions:** <https://www.doi.org/10.1016/j.jct.2019.03.011>
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- Solubility of the L-serine, L-threonine and L-isoleucine in aqueous amphiphilic micellar solutions:** <https://www.doi.org/10.1016/j.fluid.2008.05.008>
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<https://www.doi.org/10.1016/j.fluid.2007.04.004>
- Thermodynamic Properties of L-Alanine in Aqueous Solutions: Molar Heat Capacities and Apparent Molar Solvent Interactions of Homologous and Heterologous Amino Acids in Aqueous Solutions at Different Temperatures by using the Thermodynamic Methods:** <https://www.doi.org/10.1016/j.jct.2015.11.015>
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https://en.wikipedia.org/wiki/Joback_method
- Thermodynamic Properties of L-Alanine in Aqueous Solutions at Different Temperatures by using the Thermodynamic Methods:** <https://www.doi.org/10.1016/j.jct.2016.03.012>
- Thermodynamic Properties of L-Alanine in Aqueous Solutions at Different Temperatures by using the Thermodynamic Methods:** <https://www.doi.org/10.1016/j.jct.2016.03.012>
- Thermodynamic Properties of L-Alanine in Aqueous Solutions at Different Temperatures by using the Thermodynamic Methods:** <https://www.doi.org/10.1016/j.jct.2016.03.012>
- Thermodynamic Properties of L-Alanine in Aqueous Solutions at Different Temperatures by using the Thermodynamic Methods:** <https://www.doi.org/10.1016/j.jct.2016.03.012>

Legend

cpg:	Ideal gas heat capacity
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
hsubt:	Enthalpy of sublimation at a given temperature
h vap:	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
tf:	Normal melting (fusion) point
vc:	Critical Volume

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