

Glycine

Other names:	2-Aminoacetic acid Acetic acid, amino- Aciport Aminoacetic acid Aminoethanoic acid Amitone Athenon Glicoamin Gly Glycine, free base Glycine, non-medical Glycocoll Glycolixir Glycosthene Gyn-hydralin Hampshire glycine NH2CH2COOH NSC 25936 Padil
Inchi:	InChI=1S/C2H5NO2/c3-1-2(4)5/h1,3H2,(H,4,5)
InchiKey:	DHMQDGOQFOQNFH-UHFFFAOYSA-N
Formula:	C2H5NO2
SMILES:	NCC(=O)O
Mol. weight [g/mol]:	75.07
CAS:	56-40-6

Physical Properties

Property code	Value	Unit	Source
affp	883.10 ± 1.90	kJ/mol	NIST Webbook
affp	886.30 ± 3.10	kJ/mol	NIST Webbook
affp	886.50	kJ/mol	NIST Webbook
basg	851.10 ± 1.90	kJ/mol	NIST Webbook
basg	852.20	kJ/mol	NIST Webbook
basg	856.00 ± 3.00	kJ/mol	NIST Webbook
basg	855.40 ± 3.60	kJ/mol	NIST Webbook
ep	2.00 ± 6.00	J/molxK	NIST Webbook
gf	-233.33	kJ/mol	Joback Method

hf	-390.50 ± 4.60	kJ/mol	NIST Webbook
hfs	-527.50 ± 0.50	kJ/mol	NIST Webbook
hfs	-528.52 ± 0.42	kJ/mol	NIST Webbook
hfs	-537.20	kJ/mol	NIST Webbook
hfs	-528.61	kJ/mol	NIST Webbook
hfus	11.82	kJ/mol	Joback Method
hsub	138.10 ± 4.60	kJ/mol	NIST Webbook
hsub	138.10 ± 4.60	kJ/mol	NIST Webbook
hvap	54.11	kJ/mol	Joback Method
ie	9.25 ± 0.10	eV	NIST Webbook
ie	8.80	eV	NIST Webbook
ie	10.00	eV	NIST Webbook
ie	9.30	eV	NIST Webbook
ie	8.90	eV	NIST Webbook
ie	9.21 ± 0.05	eV	NIST Webbook
log10ws	0.52		Aqueous Solubility Prediction Method
logp	-0.970		Crippen Method
mcvol	56.460	ml/mol	McGowan Method
pc	6967.65	kPa	Joback Method
ss	109.20	J/molxK	NIST Webbook
ss	103.51	J/molxK	NIST Webbook
tb	463.74	K	Joback Method
tc	653.39	K	Joback Method
tf	530.49	K	Solubility of alpha-glycine in water with additives at a temperature range of (293.15 - 343.15) K: Experimental data and results of thermodynamic modeling
tf	527.85	K	Artificial neural networks as a supporting tool for compatibility study based on thermogravimetric data
vc	0.202	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	113.87	J/molxK	463.74	Joback Method
cpg	118.43	J/molxK	495.35	Joback Method
cpg	122.78	J/molxK	526.96	Joback Method
cpg	126.92	J/molxK	558.56	Joback Method

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Volumetric and Viscosity Properties of Monosaccharides in Aqueous Amino Acid Solutions at 25 and 35°C

Thermodynamic studies on interactions of a homologous series of some amino acids with trimethylamine N-oxide: Thermodynamic studies on a series of homologous α -amino acids and amino acids in some aqueous solutions of dimethyl sulfoxide and dimethyl sulfoxide-ethylene glycol mixtures

Some bio-active solutes in aqueous sodium citrate solution as stabilizing agents: Trimethylamine N-oxide, glycine and tetraammonium sulphate + NaSO₃ in beta-cyclodextrin solutions

Study of possible phase sequences of alpha-amino acids in aqueous solution by fluorescence probe in ethane-water solvent mixtures: Experimental data and thermodynamic measurement of aqueous biphasic systems composed of aminic acids in aqueous N-methylformamide, N-methylpyrrolidone and morpholine

Mixed NaCl-MgCl₂ Solutions in a Highly Concentrated Ternary System of 1-butyl-3-methylimidazolium hexafluorophosphate as a Salts Data: pharmaceutical ingredient ionic liquid (1-butyl-3-methylimidazolium amino acids 1-hexyl-3-methylimidazolium Salicylate as an Active Ingredients between some amino acids and a pharmaceutical compounds) and chemical substances of an amino acid in some ionic liquids (imidazolium) and temperatures from 298.15 K to 313.15 K

Glycine-Alcohol-NaCl-H₂O System: Volumetric and viscometric study of amino acids in aqueous sorbitol solution at different temperatures:

Infinite Dilution Binary Diffusion
Coefficients of Several α -Amino Acids
in Water over a Temperature Range
of 25 to 35 °C. Interactions of Solute
with Water Molecules for
Amino Acids in Aqueous Sodium
Bromide and Potassium Bromide
Solutions:

<https://www.doi.org/10.1021/je060149b>

<https://www.doi.org/10.1021/acs.jced.7b00647>

Legend

affp:	Proton affinity
basg:	Gas basicity
cpg:	Ideal gas heat capacity
cps:	Solid phase heat capacity
ep:	Protonation entropy at 298K
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfs:	Solid phase enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
hsub:	Enthalpy of sublimation at standard conditions
hsubt:	Enthalpy of sublimation at a given temperature
hvap:	Enthalpy of vaporization at standard conditions
ie:	Ionization energy
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
pvap:	Vapor pressure
ss:	Solid phase molar entropy at standard conditions
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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