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

InChl=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
ер	2.00 ± 6.00	J/mol×K	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	
рс	6967.65	kPa	Joback Method	
SS	109.20	J/mol×K	NIST Webbook	
SS	103.51	J/mol×K	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/mol×K	463.74	Joback Method
cpg	118.43	J/mol×K	495.35	Joback Method
cpg	122.78	J/mol×K	526.96	Joback Method
cpg	126.92	J/mol×K	558.56	Joback Method

cpg	130.87	J/mol×K	590.17	Joback Method
cpg	134.62	J/mol×K	621.78	Joback Method
cpg	138.19	J/mol×K	653.39	Joback Method
cps	95.10	J/mol×K	298.00	NIST Webbook
cps	95.00	J/mol×K	298.00	NIST Webbook
cps	99.30	J/mol×K	298.15	NIST Webbook
cps	99.20	J/mol×K	298.15	NIST Webbook
cps	100.50	J/mol×K	299.50	NIST Webbook
hsubt	137.00 ± 2.00	kJ/mol	419.50	NIST Webbook
hsubt	136.00 ± 0.40	kJ/mol	455.00	NIST Webbook
hsubt	136.40 ± 4.00	kJ/mol	462.00	NIST Webbook
hsubt	131.00 ± 2.00	kJ/mol	414.50	NIST Webbook

Correlations

Information	Value
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Property code	pvap	
Equation	$ln(Pvp) = A + B/T + C*ln(T) + D*T^2$	
Coeff. A	6.23648e+03	
Coeff. B	-3.25024e+05	
Coeff. C	-9.21946e+02	
Coeff. D	5.92592e-04	
Temperature range (K), min.	452.15	
Temperature range (K), max.	470.15	

Sources

Volumetric, compressibility and viscometric studies on sodium solvatios definition of the properties o Volumetric, compressibility and solutions at different temperatures:

https://www.doi.org/10.1016/j.tca.2015.02.014

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Bromide and Potassium Bromide
Solutions:

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Legend

affp: Proton affinity **basg:** Gas basicity

cpg: Ideal gas heat capacitycps: Solid phase heat capacityep: Protonation entropy at 298K

gf: Standard Gibbs free energy of formationhf: 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 conditionshsubt: Enthalpy of sublimation at a given temperaturehvap: 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 Pressurepvap: 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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