

dl-Serine

Other names:	(.+-.)-serine 2-amino-3-hydroxypropanoic acid DL-HOCH ₂ CH(NH ₂)COOH Serine DL-form Serine, dl- propanoic acid, 2-amino-3-hydroxy-
Inchi:	InChI=1S/C3H7NO3/c4-2(1-5)3(6)7/h2,5H,1,4H2,(H,6,7)
InchiKey:	MTCFGRXMJLQNBG-UHFFFAOYSA-N
Formula:	C ₃ H ₇ NO ₃
SMILES:	NC(CO)C(=O)O
Mol. weight [g/mol]:	105.09
CAS:	302-84-1

Physical Properties

Property code	Value	Unit	Source
chs	-1441.90 ± 1.90	kJ/mol	NIST Webbook
gf	-364.17	kJ/mol	Joback Method
hf	-493.78	kJ/mol	Joback Method
hfus	14.97	kJ/mol	Joback Method
hvap	72.63	kJ/mol	Joback Method
ie	8.70	eV	NIST Webbook
ie	10.00	eV	NIST Webbook
log10ws	1.01		Crippen Method
logp	-1.609		Crippen Method
mcvol	76.420	ml/mol	McGowan Method
pc	7014.41	kPa	Joback Method
tb	578.36	K	Joback Method
tc	760.22	K	Joback Method
tf	363.40	K	Joback Method
vc	0.271	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
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cpg	184.80	J/molxK	578.36	Joback Method
cpg	190.08	J/molxK	608.67	Joback Method
cpg	195.08	J/molxK	638.98	Joback Method
cpg	199.83	J/molxK	669.29	Joback Method
cpg	204.33	J/molxK	699.60	Joback Method
cpg	208.59	J/molxK	729.91	Joback Method
cpg	212.60	J/molxK	760.22	Joback Method
cps	132.40	J/molxK	298.15	NIST Webbook

Sources

Thermodynamics of proton dissociations from aqueous serine at temperatures from 273.15 to 315 K, Solid-Solvent Interactions of Some Amino Acids in the Presence of NaCl and DMSO: Thermodynamic Parameters and Partial Molar Volumes of Serine, Glycine, Alanine, Threonine, and Sodium Serinate: NIST Webbook.

Partial molar volumes of L-alanine, DL-serine, DL-threonine, L-histidine, Glycine, and methylglycine in water, NaCl, and DMSO aqueous solutions at effects of hydroxyl groups on binary diffusion coefficients of -amino acids in dilute aqueous solutions of serine stereoisomers: Crippen Method:

McGowan Method:

<https://www.doi.org/10.1016/j.jct.2005.07.019>

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

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

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C302841&Units=SI>

<https://www.doi.org/10.1016/j.jct.2005.03.015>

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

<https://www.doi.org/10.1016/j.fluid.2007.10.013>

<https://www.doi.org/10.1016/j.tca.2014.08.032>

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

<http://link.springer.com/article/10.1007/BF02311772>

Legend

chs:	Standard solid enthalpy of combustion
cpg:	Ideal gas heat capacity
cps:	Solid phase heat capacity
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
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
tb:	Normal Boiling Point Temperature

tc: Critical Temperature
tf: Normal melting (fusion) point
vc: Critical Volume

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