

Acetic acid, hydroxy-

Other names:	2-Hydroxyacetic acid 2-hydroxyethanoic acid ALPHA-HYDROXYACETIC ACID GLYCOLIC ACID GLYCOLLIC ACID HOCH ₂ COOH HYDROXYACETIC ACID Kyselina glykolova Kyselina hydroxyoctova NSC 166 hydroxyethanoic acid «alpha»-Hydroxyacetic acid
Inchi:	InChI=1S/C2H4O3/c3-1-2(4)5/h3H,1H2,(H,4,5)
InchiKey:	AEMRFAOFKBGASW-UHFFFAOYSA-N
Formula:	C ₂ H ₄ O ₃
SMILES:	O=C(O)CO
Mol. weight [g/mol]:	76.05
CAS:	79-14-1

Physical Properties

Property code	Value	Unit	Source
gf	-436.60	kJ/mol	Joback Method
hf	-501.65	kJ/mol	Joback Method
hfus	10.71	kJ/mol	Joback Method
hvap	60.15	kJ/mol	Joback Method
log10ws	0.98		Crippen Method
logp	-0.937		Crippen Method
mcvol	52.350	ml/mol	McGowan Method
pc	7393.34	kPa	Joback Method
tb	483.39	K	Joback Method
tc	653.36	K	Joback Method
tf	350.75	K	Thermodynamic equilibrium of hydroxyacetic acid in pure and binary solvent systems
vc	0.192	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	123.97	J/molxK	653.36	Joback Method
cpg	104.94	J/molxK	483.39	Joback Method
cpg	108.47	J/molxK	511.72	Joback Method
cpg	111.85	J/molxK	540.05	Joback Method
cpg	115.09	J/molxK	568.37	Joback Method
cpg	118.19	J/molxK	596.70	Joback Method
cpg	121.14	J/molxK	625.03	Joback Method
dvisc	0.0001395	Paxs	483.39	Joback Method
dvisc	0.0611762	Paxs	283.87	Joback Method
dvisc	0.0130420	Paxs	317.12	Joback Method
dvisc	0.0037283	Paxs	350.38	Joback Method
dvisc	0.0013242	Paxs	383.63	Joback Method
dvisc	0.0005548	Paxs	416.88	Joback Method
dvisc	0.0002643	Paxs	450.14	Joback Method
hvapt	51.80	kJ/mol	362.50	NIST Webbook

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
Isopropanol dehydration via extractive distillation using low transition temperature mixtures as entrainers:	https://www.doi.org/10.1016/j.jct.2015.02.003
KDB:	https://www.cheric.org/research/kdb/hcprop/showprop.php?cmpid=971
Thermophysical Properties and Solubility of Different Sugar-Derived Molecules in Deep Eutectic Solvents:	https://www.doi.org/10.1021/acs.jced.7b00184
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C79141&Units=SI
Low transition temperature mixtures (LTTMs) as novel entrainers in the thermodynamic equilibrium of hydroxyacetic acid in pure and binary aprotic systems:	https://www.doi.org/10.1016/j.fluid.2014.10.044
Isopropanol dehydration via Extractive Distillation Using Low Transition Temperature Mixtures as Entrainers:	https://www.doi.org/10.1016/j.jct.2017.01.004
Joback Method:	https://www.doi.org/10.1021/acs.jced.8b00228
Joback Method:	https://en.wikipedia.org/wiki/Joback_method

Legend

cp_g:	Ideal gas heat capacity
dv_{isc}:	Dynamic viscosity
g_f:	Standard Gibbs free energy of formation
h_f:	Enthalpy of formation at standard conditions
h_{fus}:	Enthalpy of fusion at standard conditions
h_{vap}:	Enthalpy of vaporization at standard conditions
h_{vapt}:	Enthalpy of vaporization at a given temperature
log₁₀w_s:	Log ₁₀ of Water solubility in mol/l
log_p:	Octanol/Water partition coefficient
mc_{vol}:	McGowan's characteristic volume
pc:	Critical Pressure
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/16-002-1/Acetic-acid-hydroxy.pdf>

Generated by Cheméo on 2024-04-25 04:44:41.121038112 +0000 UTC m=+16309530.041615424.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.