

# N-(2-Hydroxyethyl)lactamide

<b>Other names:</b>	2-Hydroxy-N-(2-hydroxyethyl)propanamide Incromectant LMEA Lactamide MEA Lactamide, N-(2-hydroxyethyl)- Lactic acid monoethanolamide Lipamide LMEA Mackamide LME Monoethanolamine lactic acid amide N-(«beta»-Hydroxyethyl)-2-hydroxypropionamide N-(«beta»-Hydroxyethyl)lactamide N-(«beta»-Hydroxyethyl)-2-hydroxypropionamide N-(«beta»-Hydroxyethyl)lactamide N-Hydroxyaethylactamid NSC 11062 Parapel LAM-100 Propanamide, 2-hydroxy-N-(2-hydroxyethyl)- Schercomid LME-100
<b>Inchi:</b>	InChI=1S/C5H11NO3/c1-4(8)5(9)6-2-3-7/h4,7-8H,2-3H2,1H3,(H,6,9)
<b>InchiKey:</b>	RZCHTMXTKQHYDT-UHFFFAOYSA-N
<b>Formula:</b>	C5H11NO3
<b>SMILES:</b>	CC(O)C(=O)NCCO
<b>Mol. weight [g/mol]:</b>	133.15
<b>CAS:</b>	5422-34-4

## Physical Properties

Property code	Value	Unit	Source
gf	-324.39	kJ/mol	Joback Method
hf	-515.38	kJ/mol	Joback Method
hfus	20.06	kJ/mol	Joback Method
hvap	72.88	kJ/mol	Joback Method
log10ws	0.48		Crippen Method
logp	-1.524		Crippen Method
mcvol	104.600	ml/mol	McGowan Method
pc	4890.21	kPa	Joback Method
tb	601.76	K	Joback Method
tc	774.06	K	Joback Method
tf	355.34	K	Joback Method

vc

0.389

m<sup>3</sup>/kmol

Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	266.13	J/mol×K	601.76	Joback Method
cpg	273.53	J/mol×K	630.48	Joback Method
cpg	280.58	J/mol×K	659.19	Joback Method
cpg	287.30	J/mol×K	687.91	Joback Method
cpg	293.68	J/mol×K	716.63	Joback Method
cpg	299.75	J/mol×K	745.35	Joback Method
cpg	305.50	J/mol×K	774.06	Joback Method

## Pressure Dependent Properties

Property code	Value	Unit	Pressure [kPa]	Source
tbrp	453.50 ± 2.50	K	0.01	NIST Webbook

## Sources

**Joback Method:**

[https://en.wikipedia.org/wiki/Joback\\_method](https://en.wikipedia.org/wiki/Joback_method)

**McGowan Method:**

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

**NIST Webbook:**

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C5422344&Units=SI>

**Crippen Method:**

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

**Crippen Method:**

[https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)

**Specific Heat Capacities of Two Functional Ionic Liquids and Two Functional Deep Eutectic Solvents for the Absorption of SO<sub>2</sub>:**

<https://www.doi.org/10.1021/acs.jced.7b00102>

## Legend

**cpg:** Ideal gas heat capacity  
**gf:** Standard Gibbs free energy of formation  
**hf:** Enthalpy of formation at standard conditions

<b>hfus:</b>	Enthalpy of fusion at standard conditions
<b>hvac:</b>	Enthalpy of vaporization at standard conditions
<b>log10ws:</b>	Log10 of Water solubility in mol/l
<b>logp:</b>	Octanol/Water partition coefficient
<b>mcvol:</b>	McGowan's characteristic volume
<b>pc:</b>	Critical Pressure
<b>tb:</b>	Normal Boiling Point Temperature
<b>tbrp:</b>	Boiling point at reduced pressure
<b>tc:</b>	Critical Temperature
<b>tf:</b>	Normal melting (fusion) point
<b>vc:</b>	Critical Volume

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