

# Acetic acid, lithium salt

Other names:	lithium acetate lithium ethanoate
Inchi:	InChI=1S/C2H4O2.Li/c1-2(3)4;/h1H3,(H,3,4);/q;+1/p-1
InchiKey:	XIXADJRWDQXREU-UHFFFAOYSA-M
Formula:	C2H3LiO2
SMILES:	CC(=O)[O-].[Li]
Mol. weight [g/mol]:	65.98
CAS:	546-89-4

## Physical Properties

Property code	Value	Unit	Source
hfus	12.60	kJ/mol	Enthalpies of formation and lattice enthalpies of alkaline metal acetates
tf	556.25	K	Lithium C1-C12 n-Alkanoates: Thermal behavior from -30C to 600C

## Sources

Salt influence on MIBK/water liquid-liquid equilibrium: Measuring influence of electrolytes on liquid-liquid equilibria: water/1-butanol and on the enthalpies of formation and lattice enthalpies of alkaline metal acetates: Lithium C1-C12 n-Alkanoates: Thermal behavior from -30C to 600C: NIST Webbook:	<a href="https://www.doi.org/10.1016/j.fluid.2015.11.018">https://www.doi.org/10.1016/j.fluid.2015.11.018</a> <a href="https://www.doi.org/10.1016/j.fluid.2016.05.001">https://www.doi.org/10.1016/j.fluid.2016.05.001</a> <a href="https://www.doi.org/10.1016/j.tca.2004.11.004">https://www.doi.org/10.1016/j.tca.2004.11.004</a> <a href="https://www.doi.org/10.1021/je400054m">https://www.doi.org/10.1021/je400054m</a> <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C546894&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C546894&amp;Units=SI</a>
Ion association and solvation behavior of some alkali metal acetates in concentrated binary and quaternary eutectic systems: Pair and triple ion structure of ethanoates in tetrahydrofuran, dimethyl sulfoxide and their binaries:	<a href="https://www.doi.org/10.1016/j.fluid.2011.05.017">https://www.doi.org/10.1016/j.fluid.2011.05.017</a> <a href="https://www.doi.org/10.1016/j.fluid.2012.02.013">https://www.doi.org/10.1016/j.fluid.2012.02.013</a>

## Legend

hfus:	Enthalpy of fusion at standard conditions
tf:	Normal melting (fusion) point

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