

# 2-Ethylbutyric acid, isohexyl ester

<b>Inchi:</b>	InChI=1S/C12H24O2/c1-5-11(6-2)12(13)14-9-7-8-10(3)4/h10-11H,5-9H2,1-4H3
<b>InchiKey:</b>	DCJUCJHQIDKEFN-UHFFFAOYSA-N
<b>Formula:</b>	C12H24O2
<b>SMILES:</b>	CCC(CC)C(=O)OCCCC(C)C
<b>Mol. weight [g/mol]:</b>	200.32

## Physical Properties

Property code	Value	Unit	Source
gf	-188.64	kJ/mol	Joback Method
hf	-546.37	kJ/mol	Joback Method
hfus	22.58	kJ/mol	Joback Method
hvap	50.69	kJ/mol	Joback Method
log10ws	-3.22		Crippen Method
logp	3.402		Crippen Method
mcvol	187.380	ml/mol	McGowan Method
pc	1877.28	kPa	Joback Method
rinpol	1252.00		NIST Webbook
rinpol	1252.00		NIST Webbook
tb	549.37	K	Joback Method
tc	725.46	K	Joback Method
tf	267.16	K	Joback Method
vc	0.720	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	464.71	J/molxK	549.37	Joback Method
cpg	481.08	J/molxK	578.72	Joback Method
cpg	496.78	J/molxK	608.07	Joback Method
cpg	511.82	J/molxK	637.42	Joback Method
cpg	526.20	J/molxK	666.76	Joback Method
cpg	539.94	J/molxK	696.11	Joback Method
cpg	553.06	J/molxK	725.46	Joback Method
dvisc	0.0062070	Paxs	267.16	Joback Method

dvisc	0.0021594	Paxs	314.19	Joback Method
dvisc	0.0009890	Paxs	361.23	Joback Method
dvisc	0.0005423	Paxs	408.26	Joback Method
dvisc	0.0003366	Paxs	455.30	Joback Method
dvisc	0.0002285	Paxs	502.33	Joback Method
dvisc	0.0001657	Paxs	549.37	Joback Method

## Sources

<b>McGowan Method:</b>	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>
<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=U340240&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U340240&amp;Units=SI</a>
<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>
<b>Crippen Method:</b>	<a href="https://www.cheméo.com/doc/models/crippen_log10ws">https://www.cheméo.com/doc/models/crippen_log10ws</a>
<b>Joback Method:</b>	<a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>

## Legend

<b>cpg:</b>	Ideal gas heat capacity
<b>dvisc:</b>	Dynamic viscosity
<b>gf:</b>	Standard Gibbs free energy of formation
<b>hf:</b>	Enthalpy of formation at standard conditions
<b>hfus:</b>	Enthalpy of fusion at standard conditions
<b>hvap:</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>rinpol:</b>	Non-polar retention indices
<b>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature
<b>tf:</b>	Normal melting (fusion) point
<b>vc:</b>	Critical Volume

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