

# Succinic acid, 3-methylbut-2-yl trans-hex-3-en-1-yl ester

Inchi:	InChI=1S/C15H26O4/c1-5-6-7-8-11-18-14(16)9-10-15(17)19-13(4)12(2)3/h6-7,12-13H,5,
InchiKey:	CYHRASGMVRBHKL-VOTSOKGWSA-N
Formula:	C15H26O4
SMILES:	CCC=CCCOC(=O)CCC(=O)OC(C)C(C)C
Mol. weight [g/mol]:	270.36

## Physical Properties

Property code	Value	Unit	Source
gf	-317.08	kJ/mol	Joback Method
hf	-735.87	kJ/mol	Joback Method
hfus	33.34	kJ/mol	Joback Method
hvap	66.48	kJ/mol	Joback Method
log10ws	-3.55		Crippen Method
logp	3.254		Crippen Method
mvol	232.790	ml/mol	McGowan Method
pc	1602.56	kPa	Joback Method
rinpol	1739.00		NIST Webbook
tb	698.46	K	Joback Method
tc	884.01	K	Joback Method
tf	368.05	K	Joback Method
vc	0.891	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	655.63	J/molxK	698.46	Joback Method
cpg	671.72	J/molxK	729.38	Joback Method
cpg	686.98	J/molxK	760.31	Joback Method
cpg	701.42	J/molxK	791.23	Joback Method
cpg	715.06	J/molxK	822.16	Joback Method
cpg	727.92	J/molxK	853.08	Joback Method
cpg	740.00	J/molxK	884.01	Joback Method
dvisc	0.0019655	Paxs	368.05	Joback Method
dvisc	0.0008087	Paxs	423.12	Joback Method

dvisc	0.0004083	Paxs	478.19	Joback Method
dvisc	0.0002374	Paxs	533.25	Joback Method
dvisc	0.0001528	Paxs	588.32	Joback Method
dvisc	0.0001060	Paxs	643.39	Joback Method
dvisc	0.0000779	Paxs	698.46	Joback Method

## Sources

<b>Joback Method:</b>	<a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>
<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=U391111&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U391111&amp;Units=SI</a>
<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</a>
<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</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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