

# Succinic acid, 4-chloro-2-methylbenzyl propyl ester

Inchi:	InChI=1S/C15H19ClO4/c1-3-8-19-14(17)6-7-15(18)20-10-12-4-5-13(16)9-11(12)2/h4-5,9
InchiKey:	FQVGKVVZNQFAZNC-UHFFFAOYSA-N
Formula:	C15H19ClO4
SMILES:	CCCOC(=O)CCC(=O)OCc1ccc(Cl)cc1C
Mol. weight [g/mol]:	298.76

## Physical Properties

Property code	Value	Unit	Source
gf	-311.20	kJ/mol	Joback Method
hf	-644.68	kJ/mol	Joback Method
hfus	37.64	kJ/mol	Joback Method
hvap	75.28	kJ/mol	Joback Method
log10ws	-4.17		Crippen Method
logp	3.425		Crippen Method
mvol	225.570	ml/mol	McGowan Method
pc	1887.08	kPa	Joback Method
rinpol	2130.00		NIST Webbook
rinpol	2130.00		NIST Webbook
tb	769.25	K	Joback Method
tc	977.26	K	Joback Method
tf	484.51	K	Joback Method
vc	0.865	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	618.97	J/molxK	769.25	Joback Method
cpg	677.58	J/molxK	942.59	Joback Method
cpg	667.71	J/molxK	907.92	Joback Method
cpg	656.93	J/molxK	873.25	Joback Method
cpg	645.21	J/molxK	838.59	Joback Method
cpg	632.56	J/molxK	803.92	Joback Method
cpg	686.52	J/molxK	977.26	Joback Method
dvisc	0.0000950	Paxs	769.25	Joback Method

dvisc	0.0001183	Paxs	721.79	Joback Method
dvisc	0.0001519	Paxs	674.34	Joback Method
dvisc	0.0002024	Paxs	626.88	Joback Method
dvisc	0.0002829	Paxs	579.42	Joback Method
dvisc	0.0004197	Paxs	531.97	Joback Method
dvisc	0.0006726	Paxs	484.51	Joback Method

## Sources

<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.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>
<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=U380878&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U380878&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>

## 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

Latest version available from:

<https://www.chemeo.com/cid/85-085-4/Succinic-acid-4-chloro-2-methylbenzyl-propyl-ester.pdf>

Generated by Cheméo on 2024-04-23 14:19:37.927606151 +0000 UTC m=+16171226.848183463.

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