

# Succinic acid, phenethyl 2-methoxyethyl ester

<b>Inchi:</b>	InChI=1S/C15H20O5/c1-18-11-12-20-15(17)8-7-14(16)19-10-9-13-5-3-2-4-6-13/h2-6H,7-
<b>InchiKey:</b>	YGZVEFQXXIPITA-UHFFFAOYSA-N
<b>Formula:</b>	C15H20O5
<b>SMILES:</b>	COCCOC(=O)CCC(=O)OCCc1ccccc1
<b>Mol. weight [g/mol]:</b>	280.32

## Physical Properties

Property code	Value	Unit	Source
gf	-385.01	kJ/mol	Joback Method
hf	-738.22	kJ/mol	Joback Method
hfus	35.41	kJ/mol	Joback Method
hvap	71.98	kJ/mol	Joback Method
log10ws	-2.02		Crippen Method
logp	1.742		Crippen Method
mcvol	219.200	ml/mol	McGowan Method
pc	1971.80	kPa	Joback Method
rinsol	2130.00		NIST Webbook
tb	744.28	K	Joback Method
tc	945.99	K	Joback Method
tf	451.78	K	Joback Method
vc	0.834	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	621.80	J/molxK	744.28	Joback Method
cpg	636.35	J/molxK	777.90	Joback Method
cpg	649.93	J/molxK	811.52	Joback Method
cpg	662.54	J/molxK	845.14	Joback Method
cpg	674.18	J/molxK	878.75	Joback Method
cpg	684.87	J/molxK	912.37	Joback Method
cpg	694.59	J/molxK	945.99	Joback Method
dvisc	0.0007872	Paxs	451.78	Joback Method
dvisc	0.0004480	Paxs	500.53	Joback Method

dvisc	0.0002818	Paxs	549.28	Joback Method
dvisc	0.0001911	Paxs	598.03	Joback Method
dvisc	0.0001375	Paxs	646.78	Joback Method
dvisc	0.0001035	Paxs	695.53	Joback Method
dvisc	0.0000809	Paxs	744.28	Joback Method

## Sources

<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.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=U358000&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U358000&amp;Units=SI</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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