

# Succinic acid, 3-iodobenzyl pentyl ester

<b>Inchi:</b>	InChI=1S/C16H21IO4/c1-2-3-4-10-20-15(18)8-9-16(19)21-12-13-6-5-7-14(17)11-13/h5-7
<b>InchiKey:</b>	GRVVXXACMPYBPK-UHFFFAOYSA-N
<b>Formula:</b>	C16H21IO4
<b>SMILES:</b>	CCCCCOC(=O)CCC(=O)OCc1cccc(I)c1
<b>Mol. weight [g/mol]:</b>	404.24

## Physical Properties

Property code	Value	Unit	Source
gf	-223.10	kJ/mol	Joback Method
hf	-561.24	kJ/mol	Joback Method
hfus	40.83	kJ/mol	Joback Method
hvap	81.83	kJ/mol	Joback Method
log10ws	-4.99		Crippen Method
logp	3.848		Crippen Method
mvol	253.240	ml/mol	McGowan Method
pc	1783.35	kPa	Joback Method
rinpol	2445.00		NIST Webbook
rinpol	2445.00		NIST Webbook
tb	842.86	K	Joback Method
tc	1064.56	K	Joback Method
tf	511.40	K	Joback Method
vc	0.960	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	698.57	J/molxK	842.86	Joback Method
cpg	754.13	J/molxK	1027.61	Joback Method
cpg	745.04	J/molxK	990.66	Joback Method
cpg	734.96	J/molxK	953.71	Joback Method
cpg	723.88	J/molxK	916.76	Joback Method
cpg	711.76	J/molxK	879.81	Joback Method
cpg	762.27	J/molxK	1064.56	Joback Method
dvisc	0.0000704	Paxs	842.86	Joback Method

dvisc	0.0000895	Paxs	787.62	Joback Method
dvisc	0.0001181	Paxs	732.37	Joback Method
dvisc	0.0001629	Paxs	677.13	Joback Method
dvisc	0.0002381	Paxs	621.89	Joback Method
dvisc	0.0003746	Paxs	566.64	Joback Method
dvisc	0.0006500	Paxs	511.40	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=U381384&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U381384&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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