

Succinic acid, 2-iodobenzyl pentyl ester

Inchi:	InChI=1S/C16H21IO4/c1-2-3-6-11-20-15(18)9-10-16(19)21-12-13-7-4-5-8-14(13)17/h4-5
InchiKey:	FMZAOADJIKXILB-UHFFFAOYSA-N
Formula:	C16H21IO4
SMILES:	CCCCCOC(=O)CCC(=O)OCc1ccccc1I
Mol. weight [g/mol]:	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
mcvol	253.240	ml/mol	McGowan Method
pc	1783.35	kPa	Joback Method
rinpol	2443.00		NIST Webbook
rinpol	2443.00		NIST Webbook
tb	842.86	K	Joback Method
tc	1064.56	K	Joback Method
tf	511.40	K	Joback Method
vc	0.960	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	698.57	J/molxK	842.86	Joback Method
cpg	711.76	J/molxK	879.81	Joback Method
cpg	723.88	J/molxK	916.76	Joback Method
cpg	734.96	J/molxK	953.71	Joback Method
cpg	745.04	J/molxK	990.66	Joback Method
cpg	754.13	J/molxK	1027.61	Joback Method
cpg	762.27	J/molxK	1064.56	Joback Method
dvisc	0.0006500	Paxs	511.40	Joback Method

dvisc	0.0003746	Paxs	566.64	Joback Method
dvisc	0.0002381	Paxs	621.89	Joback Method
dvisc	0.0001629	Paxs	677.13	Joback Method
dvisc	0.0001181	Paxs	732.37	Joback Method
dvisc	0.0000895	Paxs	787.62	Joback Method
dvisc	0.0000704	Paxs	842.86	Joback Method

Sources

Joback Method:	https://en.wikipedia.org/wiki/Joback_method
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U381103&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws

Legend

cp_g:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
h_{vap}:	Enthalpy of vaporization at standard conditions
log₁₀ws:	Log ₁₀ of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
m_{cvol}:	McGowan's characteristic volume
pc:	Critical Pressure
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/125-255-0/Succinic-acid-2-iodobenzyl-pentyl-ester.pdf>

Generated by Cheméo on 2024-04-27 14:51:41.139198775 +0000 UTC m=+16518750.059776087.

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