

Succinic acid, hexyl 2-methoxybenzyl ester

Inchi: InChI=1S/C18H26O5/c1-3-4-5-8-13-22-17(19)11-12-18(20)23-14-15-9-6-7-10-16(15)21-2
InchiKey: HRINCKLZCNWVHX-UHFFFAOYSA-N
Formula: C18H26O5
SMILES: CCCCCCOC(=O)CCC(=O)OCc1ccccc1OC
Mol. weight [g/mol]: 322.40

Physical Properties

Property code	Value	Unit	Source
gf	-369.38	kJ/mol	Joback Method
hf	-811.61	kJ/mol	Joback Method
hfus	42.79	kJ/mol	Joback Method
hvap	79.32	kJ/mol	Joback Method
log10ws	-4.38		Crippen Method
logp	3.642		Crippen Method
mcvol	261.470	ml/mol	McGowan Method
pc	1531.86	kPa	Joback Method
rinpol	2365.00		NIST Webbook
rinpol	2365.00		NIST Webbook
tb	817.90	K	Joback Method
tc	1017.91	K	Joback Method
tf	498.11	K	Joback Method
vc	1.002	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	790.00	J/molxK	817.90	Joback Method
cpg	805.25	J/molxK	851.23	Joback Method
cpg	819.39	J/molxK	884.57	Joback Method
cpg	832.44	J/molxK	917.90	Joback Method
cpg	844.40	J/molxK	951.24	Joback Method
cpg	855.27	J/molxK	984.57	Joback Method
cpg	865.06	J/molxK	1017.91	Joback Method
dvisc	0.0005083	Paxs	498.11	Joback Method

dvisc	0.0002926	Paxs	551.41	Joback Method
dvisc	0.0001857	Paxs	604.71	Joback Method
dvisc	0.0001268	Paxs	658.00	Joback Method
dvisc	0.0000917	Paxs	711.30	Joback Method
dvisc	0.0000694	Paxs	764.60	Joback Method
dvisc	0.0000544	Paxs	817.90	Joback Method

Sources

Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
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=U381209&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

Legend

cp_g:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
g_f:	Standard Gibbs free energy of formation
h_f:	Enthalpy of formation at standard conditions
h_{fus}:	Enthalpy of fusion at standard conditions
h_{vap}:	Enthalpy of vaporization at standard conditions
log₁₀w_s:	Log ₁₀ of Water solubility in mol/l
log_p:	Octanol/Water partition coefficient
mc_{vol}:	McGowan's characteristic volume
pc:	Critical Pressure
rin_{pol}:	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/89-207-4/Succinic-acid-hexyl-2-methoxybenzyl-ester.pdf>

Generated by Cheméo on 2024-04-28 07:00:00.735861977 +0000 UTC m=+16576849.656439289.

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