

Sebacic acid, phenyl undecyl ester

Inchi:	InChI=1S/C27H44O4/c1-2-3-4-5-6-7-10-13-19-24-30-26(28)22-17-11-8-9-12-18-23-27(29)
InchiKey:	CWDMSUFYEJBBQN-UHFFFAOYSA-N
Formula:	C27H44O4
SMILES:	CCCCCCCCCOC(=O)CCCCCCCC(=O)Oc1ccccc1
Mol. weight [g/mol]:	432.64

Physical Properties

Property code	Value	Unit	Source
gf	-178.97	kJ/mol	Joback Method
hf	-853.68	kJ/mol	Joback Method
hfus	65.30	kJ/mol	Joback Method
hvap	96.28	kJ/mol	Joback Method
log10ws	-8.60		Crippen Method
logp	7.787		Crippen Method
mvol	382.410	ml/mol	McGowan Method
pc	871.71	kPa	Joback Method
rinpol	3295.00		NIST Webbook
rinpol	3295.00		NIST Webbook
tb	996.42	K	Joback Method
tc	1221.91	K	Joback Method
tf	564.79	K	Joback Method
vc	1.488	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1307.01	J/molxK	996.42	Joback Method
cpg	1325.04	J/molxK	1034.00	Joback Method
cpg	1341.46	J/molxK	1071.58	Joback Method
cpg	1356.33	J/molxK	1109.17	Joback Method
cpg	1369.72	J/molxK	1146.75	Joback Method
cpg	1381.69	J/molxK	1184.33	Joback Method
cpg	1392.31	J/molxK	1221.91	Joback Method
dvisc	0.0003070	Paxs	564.79	Joback Method

dvisc	0.0001487	Paxs	636.73	Joback Method
dvisc	0.0000835	Paxs	708.67	Joback Method
dvisc	0.0000521	Paxs	780.61	Joback Method
dvisc	0.0000352	Paxs	852.54	Joback Method
dvisc	0.0000253	Paxs	924.48	Joback Method
dvisc	0.0000191	Paxs	996.42	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U354515&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Joback Method:	https://en.wikipedia.org/wiki/Joback_method

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
hvap:	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

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