

Sebacic acid, oct-3-yl propyl ester

Inchi:	InChI=1S/C21H40O4/c1-4-7-12-15-19(6-3)25-21(23)17-14-11-9-8-10-13-16-20(22)24-18
InchiKey:	DDLBEERQXVUQQG-UHFFFAOYSA-N
Formula:	C21H40O4
SMILES:	CCCCC(CC)OC(=O)CCCCCCCCC(=O)OCCC
Mol. weight [g/mol]:	356.54

Physical Properties

Property code	Value	Unit	Source
gf	-344.34	kJ/mol	Joback Method
hf	-971.65	kJ/mol	Joback Method
hfus	52.20	kJ/mol	Joback Method
hvap	80.26	kJ/mol	Joback Method
log10ws	-6.45		Crippen Method
logp	5.963		Crippen Method
mcvol	321.630	ml/mol	McGowan Method
pc	1018.13	kPa	Joback Method
rinpol	2295.00		NIST Webbook
rinpol	2295.00		NIST Webbook
tb	832.02	K	Joback Method
tc	1019.80	K	Joback Method
tf	455.75	K	Joback Method
vc	1.254	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1030.73	J/mol×K	832.02	Joback Method
cpg	1049.46	J/mol×K	863.32	Joback Method
cpg	1067.07	J/mol×K	894.61	Joback Method
cpg	1083.56	J/mol×K	925.91	Joback Method
cpg	1098.97	J/mol×K	957.21	Joback Method
cpg	1113.31	J/mol×K	988.51	Joback Method
cpg	1126.61	J/mol×K	1019.80	Joback Method
dvisc	0.0008818	Pa×s	455.75	Joback Method

dvisc	0.0003912	Pa×s	518.46	Joback Method
dvisc	0.0002068	Pa×s	581.17	Joback Method
dvisc	0.0001238	Pa×s	643.88	Joback Method
dvisc	0.0000812	Pa×s	706.60	Joback Method
dvisc	0.0000570	Pa×s	769.31	Joback Method
dvisc	0.0000422	Pa×s	832.02	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U416826&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Joback Method:	https://en.wikipedia.org/wiki/Joback_method

Legend

cpg:	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:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	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/80-170-4/Sebacic-acid-oct-3-yl-propyl-ester.pdf>

Generated by Cheméo on 2025-12-23 09:41:19.594134393 +0000 UTC m=+6231077.124175046.

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