

Glutaric acid, 2-methylhex-3-yl 2-propylpentyl ester

Inchi:	InChI=1S/C20H38O4/c1-6-10-17(11-7-2)15-23-19(21)13-9-14-20(22)24-18(12-8-3)16(4)5
InchiKey:	IWGMBUQQZRACEC-UHFFFAOYSA-N
Formula:	C20H38O4
SMILES:	CCCC(CCC)COC(=O)CCCC(=O)OC(CCC)C(C)C
Mol. weight [g/mol]:	342.51

Physical Properties

Property code	Value	Unit	Source
gf	-357.64	kJ/mol	Joback Method
hf	-961.57	kJ/mol	Joback Method
hfus	42.56	kJ/mol	Joback Method
hvap	77.26	kJ/mol	Joback Method
log10ws	-5.55		Crippen Method
logp	5.284		Crippen Method
mcvol	307.540	ml/mol	McGowan Method
pc	1094.99	kPa	Joback Method
rinpol	2176.00		NIST Webbook
tb	808.26	K	Joback Method
tc	994.56	K	Joback Method
tf	414.48	K	Joback Method
vc	1.185	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	970.62	J/molxK	808.26	Joback Method
cpg	989.13	J/molxK	839.31	Joback Method
cpg	1006.54	J/molxK	870.36	Joback Method
cpg	1022.88	J/molxK	901.41	Joback Method
cpg	1038.15	J/molxK	932.46	Joback Method
cpg	1052.38	J/molxK	963.51	Joback Method
cpg	1065.59	J/molxK	994.56	Joback Method
dvisc	0.0014761	Paxs	414.48	Joback Method
dvisc	0.0005410	Paxs	480.11	Joback Method

dvisc	0.0002524	Paxs	545.74	Joback Method
dvisc	0.0001387	Paxs	611.37	Joback Method
dvisc	0.0000856	Paxs	677.00	Joback Method
dvisc	0.0000575	Paxs	742.63	Joback Method
dvisc	0.0000412	Paxs	808.26	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U377134&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
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772

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/35-236-1/Glutaric-acid-2-methylhex-3-yl-2-propylpentyl-ester.pdf>

Generated by Cheméo on 2024-04-20 03:05:45.931276466 +0000 UTC m=+15871594.851853782.

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