

Glutaric acid, 2-norbornyl 2-ethylhexyl ester

Inchi: InChI=1S/C20H34O4/c1-3-5-7-15(4-2)14-23-19(21)8-6-9-20(22)24-18-13-16-10-11-17(18)
InchiKey: CNRTYRMHGVYABM-UHFFFAOYSA-N
Formula: C20H34O4
SMILES: CCCCC(CC)COC(=O)CCCC(=O)OC1CC2CCC1C2
Mol. weight [g/mol]: 338.48

Physical Properties

Property code	Value	Unit	Source
gf	-251.07	kJ/mol	Joback Method
hf	-831.91	kJ/mol	Joback Method
hfus	44.85	kJ/mol	Joback Method
hvap	77.73	kJ/mol	Joback Method
log10ws	-5.10		Crippen Method
logp	4.648		Crippen Method
mvol	285.820	ml/mol	McGowan Method
pc	1288.36	kPa	Joback Method
rinpol	2342.00		NIST Webbook
rinpol	2342.00		NIST Webbook
tb	822.22	K	Joback Method
tc	1019.23	K	Joback Method
tf	472.60	K	Joback Method
vc	1.103	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	945.56	J/molxK	822.22	Joback Method
cpg	1028.88	J/molxK	986.40	Joback Method
cpg	1014.31	J/molxK	953.56	Joback Method
cpg	998.75	J/molxK	920.73	Joback Method
cpg	982.13	J/molxK	887.89	Joback Method
cpg	964.42	J/molxK	855.06	Joback Method
cpg	1042.50	J/molxK	1019.23	Joback Method
dvisc	0.0004023	Paxs	822.22	Joback Method

dvisc	0.0004800	Paxs	763.95	Joback Method
dvisc	0.0005895	Paxs	705.68	Joback Method
dvisc	0.0007514	Paxs	647.41	Joback Method
dvisc	0.0010047	Paxs	589.14	Joback Method
dvisc	0.0014320	Paxs	530.87	Joback Method
dvisc	0.0022274	Paxs	472.60	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=U405493&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
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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