

Glutaric acid, 2-methylpent-3-yl dodec-9-yn-1-yl ester

Inchi:	InChI=1S/C23H40O4/c1-5-7-8-9-10-11-12-13-14-15-19-26-22(24)17-16-18-23(25)27-21(
InchiKey:	SPDUNZFXTSTKQW-UHFFFAOYSA-N
Formula:	C23H40O4
SMILES:	CCC#CCCCCCCCOC(=O)CCCC(=O)OC(CC)C(C)C
Mol. weight [g/mol]:	380.56

Physical Properties

Property code	Value	Unit	Source
gf	-127.14	kJ/mol	Joback Method
hf	-745.91	kJ/mol	Joback Method
hfus	56.98	kJ/mol	Joback Method
hvap	86.48	kJ/mol	Joback Method
log10ws	-6.84		Crippen Method
logp	5.822		Crippen Method
mvol	341.210	ml/mol	McGowan Method
pc	1002.71	kPa	Joback Method
rinpol	2589.00		NIST Webbook
rinpol	2589.00		NIST Webbook
tb	886.34	K	Joback Method
tc	1086.67	K	Joback Method
tf	569.39	K	Joback Method
vc	1.321	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1103.44	J/molxK	886.34	Joback Method
cpg	1121.84	J/molxK	919.73	Joback Method
cpg	1138.97	J/molxK	953.12	Joback Method
cpg	1154.84	J/molxK	986.50	Joback Method
cpg	1169.50	J/molxK	1019.89	Joback Method
cpg	1182.95	J/molxK	1053.28	Joback Method
cpg	1195.24	J/molxK	1086.67	Joback Method

Sources

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
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U393941&Units=SI

Legend

cpg:	Ideal gas heat capacity
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
rinpola:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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