

Glycerol, 2,3-dimethyl, 1-(15-methylhexadecanoate)

Inchi:	InChI=1S/C22H44O4/c1-20(2)16-14-12-10-8-6-5-7-9-11-13-15-17-22(23)26-19-21(25-4)
InchiKey:	YNJFCKMONSQIOI-UHFFFAOYSA-N
Formula:	C22H44O4
SMILES:	COCC(COC(=O)CCCCCCCCCCCCC(C)C)OC
Mol. weight [g/mol]:	372.58

Physical Properties

Property code	Value	Unit	Source
gf	-314.44	kJ/mol	Joback Method
hf	-1017.21	kJ/mol	Joback Method
hfus	50.85	kJ/mol	Joback Method
hvap	77.77	kJ/mol	Joback Method
log10ws	-5.94		Crippen Method
logp	5.918		Crippen Method
mcvol	340.020	ml/mol	McGowan Method
pc	917.16	kPa	Joback Method
rinpol	2334.00		NIST Webbook
rinpol	2334.00		NIST Webbook
tb	823.01	K	Joback Method
tc	1008.16	K	Joback Method
tf	424.32	K	Joback Method
vc	1.315	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1099.11	J/molxK	823.01	Joback Method
cpg	1189.15	J/molxK	977.30	Joback Method
cpg	1173.50	J/molxK	946.44	Joback Method
cpg	1156.68	J/molxK	915.59	Joback Method
cpg	1138.69	J/molxK	884.73	Joback Method
cpg	1119.50	J/molxK	853.87	Joback Method
cpg	1203.65	J/molxK	1008.16	Joback Method
dvisc	0.0000263	Paxs	823.01	Joback Method

dvisc	0.0000365	Paxs	756.56	Joback Method
dvisc	0.0000539	Paxs	690.11	Joback Method
dvisc	0.0000866	Paxs	623.66	Joback Method
dvisc	0.0001558	Paxs	557.22	Joback Method
dvisc	0.0003285	Paxs	490.77	Joback Method
dvisc	0.0008750	Paxs	424.32	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=R56450&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
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772

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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