

Isovaleric acid, eicosyl ester

Inchi:	InChI=1S/C25H50O2/c1-4-5-6-7-8-9-10-11-12-13-14-15-16-17-18-19-20-21-22-27-25(26
InchiKey:	LPGKUHHQQXQXFT-UHFFFAOYSA-N
Formula:	C25H50O2
SMILES:	CCCCCCCCCCCCCCCCCCCCOC(=O)CC(C)C
Mol. weight [g/mol]:	382.66
CAS:	202659-51-6

Physical Properties

Property code	Value	Unit	Source
gf	-76.74	kJ/mol	Joback Method
hf	-809.41	kJ/mol	Joback Method
hfus	59.77	kJ/mol	Joback Method
hvap	80.01	kJ/mol	Joback Method
log10ws	-8.91		Crippen Method
logp	8.617		Crippen Method
mcvol	370.550	ml/mol	McGowan Method
pc	787.27	kPa	Joback Method
rinpol	2643.30		NIST Webbook
rinpol	2643.30		NIST Webbook
tb	847.25	K	Joback Method
tc	1037.37	K	Joback Method
tf	428.67	K	Joback Method
vc	1.454	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1221.19	J/mol×K	847.25	Joback Method
cpg	1243.42	J/mol×K	878.94	Joback Method
cpg	1264.39	J/mol×K	910.62	Joback Method
cpg	1284.15	J/mol×K	942.31	Joback Method
cpg	1302.75	J/mol×K	974.00	Joback Method
cpg	1320.21	J/mol×K	1005.69	Joback Method
cpg	1336.60	J/mol×K	1037.37	Joback Method

dvisc	0.0011387	Paxs	428.67	Joback Method
dvisc	0.0004178	Paxs	498.43	Joback Method
dvisc	0.0001961	Paxs	568.20	Joback Method
dvisc	0.0001086	Paxs	637.96	Joback Method
dvisc	0.0000676	Paxs	707.72	Joback Method
dvisc	0.0000458	Paxs	777.49	Joback Method
dvisc	0.0000331	Paxs	847.25	Joback Method

Sources

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
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C202659516&Units=SI

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
mvol:	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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