Pimelic acid, 2-(2-methoxyethyl)heptyl pentyl ester

Inchi: InChI=1S/C22H42O5/c1-4-6-9-13-20(16-18-25-3)19-27-22(24)15-11-8-10-14-21(23)26-1

InchiKey: KAKYHODGZDUMIC-UHFFFAOYSA-N

Formula: C22H42O5

SMILES: CCCCCOC(=O)CCCCCC(=O)OCC(CCCCC)CCOC

Mol. weight [g/mol]: 386.57

Physical Properties

Property code	Value	Unit	Source
gf	-440.92	kJ/mol	Joback Method
hf	-1124.51	kJ/mol	Joback Method
hfus	55.97	kJ/mol	Joback Method
hvap	84.90	kJ/mol	Joback Method
log10ws	-5.60		Crippen Method
logp	5.447		Crippen Method
mcvol	341.590	ml/mol	McGowan Method
pc	948.50	kPa	Joback Method
rinpol	2551.00		NIST Webbook
rinpol	2551.00		NIST Webbook
tb	877.32	K	Joback Method
tc	1074.12	K	Joback Method
tf	489.25	K	Joback Method
vc	1.327	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1124.56	J/mol×K	877.32	Joback Method
cpg	1205.79	J/mol×K	1041.32	Joback Method
cpg	1192.16	J/mol×K	1008.52	Joback Method
cpg	1177.23	J/mol×K	975.72	Joback Method
cpg	1161.00	J/mol×K	942.92	Joback Method
cpg	1143.45	J/mol×K	910.12	Joback Method
cpg	1218.14	J/mol×K	1074.12	Joback Method
dvisc	0.0000270	Pa×s	877.32	Joback Method

dvisc	0.0000363	Pa×s	812.64	Joback Method
dvisc	0.0000515	Pa×s	747.96	Joback Method
dvisc	0.0000780	Pa×s	683.28	Joback Method
dvisc	0.0001288	Pa×s	618.61	Joback Method
dvisc	0.0002392	Pa×s	553.93	Joback Method
dvisc	0.0005231	Pa×s	489.25	Joback Method

Sources

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=U406734&Units=SI
Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l

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