

2-(2-[2-(Propionyloxy)ethoxy]ethoxy)ethyl propionate

Other names:	Triethylene glycol, dipropionate
Inchi:	InChI=1S/C12H22O6/c1-3-11(13)17-9-7-15-5-6-16-8-10-18-12(14)4-2/h3-10H2,1-2H3
InchiKey:	AWKXKNCCQLNZDB-UHFFFAOYSA-N
Formula:	C12H22O6
SMILES:	CCC(=O)OCCOCCOCCOC(=O)CC
Mol. weight [g/mol]:	262.30
CAS:	141-34-4

Physical Properties

Property code	Value	Unit	Source
gf	-627.68	kJ/mol	Joback Method
hf	-1045.05	kJ/mol	Joback Method
hfus	34.79	kJ/mol	Joback Method
hvap	65.44	kJ/mol	Joback Method
log10ws	-0.75		Crippen Method
logp	0.926		Crippen Method
mcvol	206.560	ml/mol	McGowan Method
pc	1870.79	kPa	Joback Method
tb	671.38	K	Joback Method
tc	848.52	K	Joback Method
tf	413.78	K	Joback Method
vc	0.791	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	570.81	J/molxK	671.38	Joback Method
cpg	585.08	J/molxK	700.90	Joback Method
cpg	598.70	J/molxK	730.43	Joback Method
cpg	611.66	J/molxK	759.95	Joback Method
cpg	623.94	J/molxK	789.47	Joback Method
cpg	635.53	J/molxK	819.00	Joback Method
cpg	646.41	J/molxK	848.52	Joback Method
dvisc	0.0008628	Paxs	413.78	Joback Method

dvisc	0.0004978	Paxs	456.71	Joback Method
dvisc	0.0003156	Paxs	499.65	Joback Method
dvisc	0.0002151	Paxs	542.58	Joback Method
dvisc	0.0001551	Paxs	585.51	Joback Method
dvisc	0.0001169	Paxs	628.45	Joback Method
dvisc	0.0000914	Paxs	671.38	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=C141344&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
g_f:	Standard Gibbs free energy of formation
h_f:	Enthalpy of formation at standard conditions
h_{fus}:	Enthalpy of fusion at standard conditions
h_{vap}:	Enthalpy of vaporization at standard conditions
log₁₀ws:	Log ₁₀ of Water solubility in mol/l
log_p:	Octanol/Water partition coefficient
mc_{vol}:	McGowan's characteristic volume
pc:	Critical Pressure
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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