

Diglycolic acid, 2-pentyl undecyl ester

Inchi:

InchiKey:

Formula:

SMILES:

Mol. weight [g/mol]:

InChI=1S/C20H38O5/c1-4-6-7-8-9-10-11-12-13-15-24-19(21)16-23-17-20(22)25-18(3)14

BKCPPCXCXCFHIXAQ-UHFFFAOYSA-N

C20H38O5

CCCCCCCCCCCCOC(=O)COCC(=O)OC(C)CCC

358.51

Physical Properties

Property code	Value	Unit	Source
gf	-457.76	kJ/mol	Joback Method
hf	-1083.23	kJ/mol	Joback Method
hfus	50.80	kJ/mol	Joback Method
hvap	80.45	kJ/mol	Joback Method
log10ws	-5.12		Crippen Method
logp	4.809		Crippen Method
mcvol	313.410	ml/mol	McGowan Method
pc	1071.46	kPa	Joback Method
rinpol	2905.00		NIST Webbook
rinpol	2905.00		NIST Webbook
tb	831.56	K	Joback Method
tc	1019.41	K	Joback Method
tf	466.71	K	Joback Method
vc	1.216	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1001.32	J/mol×K	831.56	Joback Method
cpg	1019.40	J/mol×K	862.87	Joback Method
cpg	1036.34	J/mol×K	894.18	Joback Method
cpg	1052.14	J/mol×K	925.48	Joback Method
cpg	1066.81	J/mol×K	956.79	Joback Method
cpg	1080.36	J/mol×K	988.10	Joback Method
cpg	1092.80	J/mol×K	1019.41	Joback Method
dvisc	0.0006668	Pa×s	466.71	Joback Method

dvisc	0.0003109	Pa×s	527.52	Joback Method
dvisc	0.0001697	Pa×s	588.33	Joback Method
dvisc	0.0001037	Pa×s	649.13	Joback Method
dvisc	0.0000690	Pa×s	709.94	Joback Method
dvisc	0.0000489	Pa×s	770.75	Joback Method
dvisc	0.0000365	Pa×s	831.56	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U382342&Units=SI
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

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
log_{10ws}:	Log10 of Water solubility in mol/l
log_p:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
p_c:	Critical Pressure
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/87-588-4/Diglycolic-acid-2-pentyl-undecyl-ester.pdf>

Generated by Cheméo on 2025-12-05 16:41:19.609403244 +0000 UTC m=+4701077.139443944.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.