

Benzoic acid, 3-propyloxy-, propyl ester

Inchi:

lnchl=1S/C13H18O3/c1-3-8-15-12-7-5-6-11(10-12)13(14)16-9-4-2/h5-7,10H,3-4,8-9H2,1

InchiKey:

SXQYCRYTLYTBBS-UHFFFAOYSA-N

Formula:

C13H18O3

SMILES:

CCCOc(=O)c1cccc(OCCC)c1

Mol. weight [g/mol]:

222.28

Physical Properties

Property code	Value	Unit	Source
gf	-177.56	kJ/mol	Joback Method
hf	-463.61	kJ/mol	Joback Method
hfus	27.05	kJ/mol	Joback Method
hvap	59.04	kJ/mol	Joback Method
log10ws	-3.51		Crippen Method
logp	3.042		Crippen Method
mcvol	183.580	ml/mol	McGowan Method
pc	2222.89	kPa	Joback Method
rinpol	1685.00		NIST Webbook
tb	627.21	K	Joback Method
tc	829.16	K	Joback Method
tf	369.60	K	Joback Method
vc	0.698	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	469.42	J/mol×K	627.21	Joback Method
cpg	484.69	J/mol×K	660.87	Joback Method
cpg	499.15	J/mol×K	694.53	Joback Method
cpg	512.81	J/mol×K	728.19	Joback Method
cpg	525.67	J/mol×K	761.85	Joback Method
cpg	537.75	J/mol×K	795.50	Joback Method
cpg	549.04	J/mol×K	829.16	Joback Method
dvisc	0.0012215	Pa×s	369.60	Joback Method
dvisc	0.0006941	Pa×s	412.54	Joback Method

dvisc	0.0004388	Pa×s	455.47	Joback Method
dvisc	0.0003002	Pa×s	498.41	Joback Method
dvisc	0.0002181	Pa×s	541.34	Joback Method
dvisc	0.0001661	Pa×s	584.28	Joback Method
dvisc	0.0001313	Pa×s	627.21	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U374556&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

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

Latest version available from:

<https://www.chemeo.com/cid/42-496-5/Benzoic-acid-3-propyloxy-propyl-ester.pdf>

Generated by Cheméo on 2025-12-22 19:35:03.849241363 +0000 UTC m=+6180301.379282026.

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