

Fumaric acid, 2-butyl pentyl ester

Inchi:

InchiKey:

Formula:

SMILES:

Mol. weight [g/mol]:

InChI=1S/C13H22O4/c1-4-6-7-10-16-12(14)8-9-13(15)17-11(3)5-2/h8-9,11H,4-7,10H2,1-

FKYWKWVYRJZSPT-CMDGGGOBGSA-N

C13H22O4

CCCCCOC(=O)C=CC(=O)OC(C)CC

242.31

Physical Properties

Property code	Value	Unit	Source
gf	-331.48	kJ/mol	Joback Method
hf	-689.31	kJ/mol	Joback Method
hfus	31.68	kJ/mol	Joback Method
hvap	62.41	kJ/mol	Joback Method
log10ws	-2.95		Crippen Method
logp	2.618		Crippen Method
mcvol	204.610	ml/mol	McGowan Method
pc	1867.55	kPa	Joback Method
rinpol	1617.00		NIST Webbook
rinpol	1617.00		NIST Webbook
tb	653.14	K	Joback Method
tc	838.21	K	Joback Method
tf	360.51	K	Joback Method
vc	0.785	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	547.34	J/mol×K	653.14	Joback Method
cpg	614.70	J/mol×K	807.36	Joback Method
cpg	602.64	J/mol×K	776.52	Joback Method
cpg	589.89	J/mol×K	745.67	Joback Method
cpg	576.43	J/mol×K	714.83	Joback Method
cpg	562.25	J/mol×K	683.98	Joback Method
cpg	626.08	J/mol×K	838.21	Joback Method
dvisc	0.0001079	Pa×s	653.14	Joback Method

dvisc	0.0001431	Pa×s	604.37	Joback Method
dvisc	0.0001995	Pa×s	555.60	Joback Method
dvisc	0.0002963	Pa×s	506.82	Joback Method
dvisc	0.0004789	Pa×s	458.05	Joback Method
dvisc	0.0008679	Pa×s	409.28	Joback Method
dvisc	0.0018471	Pa×s	360.51	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=U348641&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws

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

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