2-Butenedioic acid (Z)-, dipropyl ester

Other names: dipropyl maleate
Inchi: InChI=1S/C10H16O4/c1-3-7-13-9(11)5-6-10(12)14-8-4-2/h5-6H,3-4,7-8H2,1-2H3/b6-5-
InchiKey: DSTWFRCNXMNXTR-WAYWQWQTSA-N
Formula: C10H16O4
SMILES: CCCOC(=O)C=CC(=O)OCCC
Mol. weight [g/mol]: 200.23
CAS: 2432-63-5

Physical Properties

Property code	Value	Unit	Source
gf	-354.30	kJ/mol	Joback Method
hf	-622.11	kJ/mol	Joback Method
hfus	27.43	kJ/mol	Joback Method
hvap	56.12	kJ/mol	Joback Method
log10ws	-1.59		Crippen Method
logp	1.449		Crippen Method
mcvol	162.340	ml/mol	McGowan Method
pc	2417.12	kPa	Joback Method
rinpol	1360.00		NIST Webbook
rinpol	1360.00		NIST Webbook
tb	584.94	K	Joback Method
tc	772.08	K	Joback Method
tf	341.70	K	Joback Method
vc	0.624	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	395.88	J/mol×K	584.94	Joback Method
cpg	408.59	J/mol×K	616.13	Joback Method
cpg	420.72	J/mol×K	647.32	Joback Method
cpg	432.29	J/mol×K	678.51	Joback Method
cpg	443.29	J/mol×K	709.70	Joback Method
cpg	453.74	J/mol×K	740.89	Joback Method

cpg	463.64	J/mol×K	772.08	Joback Method
dvisc	0.0017602	Pa×s	341.70	Joback Method
dvisc	0.0009565	Pa×s	382.24	Joback Method
dvisc	0.0005843	Pa×s	422.78	Joback Method
dvisc	0.0003891	Pa×s	463.32	Joback Method
dvisc	0.0002766	Pa×s	503.86	Joback Method
dvisc	0.0002069	Pa×s	544.40	Joback Method
dvisc	0.0001611	Pa×s	584.94	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=C2432635&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

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