

Glutaric acid, butyl cis-hex-3-enyl ester

Inchi:	InChI=1S/C15H26O4/c1-3-5-7-8-13-19-15(17)11-9-10-14(16)18-12-6-4-2/h5,7H,3-4,6,8-
InchiKey:	YAZJWXLXRBNZES-ALCCZGGFSA-N
Formula:	C15H26O4
SMILES:	CCC=CCCOC(=O)CCCC(=O)OCCCC
Mol. weight [g/mol]:	270.36

Physical Properties

Property code	Value	Unit	Source
gf	-312.20	kJ/mol	Joback Method
hf	-725.31	kJ/mol	Joback Method
hfus	40.38	kJ/mol	Joback Method
hvap	67.25	kJ/mol	Joback Method
log10ws	-3.68		Crippen Method
logp	3.399		Crippen Method
mvol	232.790	ml/mol	McGowan Method
pc	1582.23	kPa	Joback Method
rinpol	1899.00		NIST Webbook
rinpol	1899.00		NIST Webbook
tb	699.34	K	Joback Method
tc	880.04	K	Joback Method
tf	398.05	K	Joback Method
vc	0.903	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	654.66	J/molxK	699.34	Joback Method
cpg	670.28	J/molxK	729.46	Joback Method
cpg	685.13	J/molxK	759.57	Joback Method
cpg	699.22	J/molxK	789.69	Joback Method
cpg	712.57	J/molxK	819.81	Joback Method
cpg	725.18	J/molxK	849.93	Joback Method
cpg	737.08	J/molxK	880.04	Joback Method
dvisc	0.0012706	Paxs	398.05	Joback Method

dvisc	0.0006402	Paxs	448.26	Joback Method
dvisc	0.0003703	Paxs	498.48	Joback Method
dvisc	0.0002368	Paxs	548.69	Joback Method
dvisc	0.0001632	Paxs	598.91	Joback Method
dvisc	0.0001192	Paxs	649.12	Joback Method
dvisc	0.0000910	Paxs	699.34	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=U359964&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
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:	Log ₁₀ of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
m_{cvol}:	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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