

Glutaric acid, 2-ethylhexyl but-3-en-1-yl ester

Inchi:	InChI=1S/C17H30O4/c1-4-7-10-15(6-3)14-21-17(19)12-9-11-16(18)20-13-8-5-2/h5,15H,2
InchiKey:	GVOGZJSGLRGSED-UHFFFAOYSA-N
Formula:	C17H30O4
SMILES:	C=CCCOC(=O)CCCC(=O)OCC(CC)CCCC
Mol. weight [g/mol]:	298.42

Physical Properties

Property code	Value	Unit	Source
gf	-290.18	kJ/mol	Joback Method
hf	-763.66	kJ/mol	Joback Method
hfus	40.56	kJ/mol	Joback Method
hvap	70.69	kJ/mol	Joback Method
log10ws	-4.28		Crippen Method
logp	4.036		Crippen Method
mvol	260.970	ml/mol	McGowan Method
pc	1363.65	kPa	Joback Method
rinpol	1837.00		NIST Webbook
rinpol	1837.00		NIST Webbook
tb	737.18	K	Joback Method
tc	917.39	K	Joback Method
tf	408.91	K	Joback Method
vc	1.010	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	766.33	J/molxK	737.18	Joback Method
cpg	840.98	J/molxK	887.36	Joback Method
cpg	827.74	J/molxK	857.32	Joback Method
cpg	813.66	J/molxK	827.29	Joback Method
cpg	798.75	J/molxK	797.25	Joback Method
cpg	782.97	J/molxK	767.22	Joback Method
cpg	853.40	J/molxK	917.39	Joback Method
dvisc	0.0000793	Paxs	737.18	Joback Method

dvisc	0.0001054	Paxs	682.47	Joback Method
dvisc	0.0001473	Paxs	627.76	Joback Method
dvisc	0.0002194	Paxs	573.04	Joback Method
dvisc	0.0003555	Paxs	518.33	Joback Method
dvisc	0.0006455	Paxs	463.62	Joback Method
dvisc	0.0013750	Paxs	408.91	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=U394038&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071

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
log₁₀w_s:	Log ₁₀ of Water solubility in mol/l
log_p:	Octanol/Water partition coefficient
mc_{vol}:	McGowan's characteristic volume
pc:	Critical Pressure
rin_{pol}:	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/90-253-1/Glutaric-acid-2-ethylhexyl-but-3-en-1-yl-ester.pdf>

Generated by Cheméo on 2024-04-26 02:25:11.338160178 +0000 UTC m=+16387560.258737492.

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