

2-Ethylbutyric acid, 2-naphthylmethyl ester

Inchi:	InChI=1S/C17H20O2/c1-3-14(4-2)17(18)19-12-13-9-10-15-7-5-6-8-16(15)11-13/h5-11,14
InchiKey:	FNCSRCURYMNTDQ-UHFFFAOYSA-N
Formula:	C17H20O2
SMILES:	CCC(CC)C(=O)OCc1ccc2ccccc2c1
Mol. weight [g/mol]:	256.34

Physical Properties

Property code	Value	Unit	Source
gf	65.33	kJ/mol	Joback Method
hf	-228.16	kJ/mol	Joback Method
hfus	29.72	kJ/mol	Joback Method
hvap	66.78	kJ/mol	Joback Method
log10ws	-5.29		Crippen Method
logp	4.319		Crippen Method
mcvol	214.610	ml/mol	McGowan Method
pc	2001.91	kPa	Joback Method
rinpola	2035.00		NIST Webbook
tb	714.85	K	Joback Method
tc	933.61	K	Joback Method
tf	410.15	K	Joback Method
vc	0.820	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	591.06	J/molxK	714.85	Joback Method
cpg	607.25	J/molxK	751.31	Joback Method
cpg	622.36	J/molxK	787.77	Joback Method
cpg	636.46	J/molxK	824.23	Joback Method
cpg	649.60	J/molxK	860.69	Joback Method
cpg	661.85	J/molxK	897.15	Joback Method
cpg	673.26	J/molxK	933.61	Joback Method
dvisc	0.0014493	Paxs	410.15	Joback Method
dvisc	0.0008465	Paxs	460.93	Joback Method

dvisc	0.0005501	Paxs	511.72	Joback Method
dvisc	0.0003864	Paxs	562.50	Joback Method
dvisc	0.0002878	Paxs	613.28	Joback Method
dvisc	0.0002242	Paxs	664.07	Joback Method
dvisc	0.0001810	Paxs	714.85	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=U370561&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

Latest version available from:

<https://www.chemeo.com/cid/22-268-0/2-Ethylbutyric-acid-2-naphthylmethyl-ester.pdf>

Generated by Cheméo on 2024-05-08 21:01:49.442649211 +0000 UTC m=+17491358.363226522.

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