

Pseudoisoeugenyl 2-ethylbutyrate I

Inchi:	InChI=1S/C16H22O3/c1-5-8-12-9-10-14(15(11-12)18-4)19-16(17)13(6-2)7-3/h5,8-11,13H
InchiKey:	KEANZVAQMYREAC-UHFFFAOYSA-N
Formula:	C16H22O3
SMILES:	CC=Cc1ccc(OC(=O)C(CC)CC)c(OC)c1
Mol. weight [g/mol]:	262.34

Physical Properties

Property code	Value	Unit	Source
gf	-84.15	kJ/mol	Joback Method
hf	-425.06	kJ/mol	Joback Method
hfus	31.11	kJ/mol	Joback Method
hvap	65.95	kJ/mol	Joback Method
log10ws	-4.58		Crippen Method
logp	4.070		Crippen Method
mvol	221.550	ml/mol	McGowan Method
pc	1789.39	kPa	Joback Method
rinpol	1736.00		NIST Webbook
rinpol	1736.00		NIST Webbook
tb	704.55	K	Joback Method
tc	910.32	K	Joback Method
tf	395.85	K	Joback Method
vc	0.840	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	605.42	J/molxK	704.55	Joback Method
cpg	621.74	J/molxK	738.85	Joback Method
cpg	637.10	J/molxK	773.14	Joback Method
cpg	651.53	J/molxK	807.44	Joback Method
cpg	665.05	J/molxK	841.73	Joback Method
cpg	677.67	J/molxK	876.03	Joback Method
cpg	689.42	J/molxK	910.32	Joback Method
dvisc	0.0009069	Paxs	395.85	Joback Method

dvisc	0.0004713	Paxs	447.30	Joback Method
dvisc	0.0002804	Paxs	498.75	Joback Method
dvisc	0.0001838	Paxs	550.20	Joback Method
dvisc	0.0001295	Paxs	601.65	Joback Method
dvisc	0.0000964	Paxs	653.10	Joback Method
dvisc	0.0000750	Paxs	704.55	Joback Method

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
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=R510766&Units=SI

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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