

Cyclopropanecarboxylic acid, trans-2-phenyl-, octyl ester

Inchi:	InChI=1S/C18H26O2/c1-2-3-4-5-6-10-13-20-18(19)17-14-16(17)15-11-8-7-9-12-15/h7-9,
InchiKey:	ANWXAQWYWWNGX-UHFFFAOYSA-N
Formula:	C18H26O2
SMILES:	CCCCCCCCOC(=O)C1CC1c1ccccc1
Mol. weight [g/mol]:	274.40

Physical Properties

Property code	Value	Unit	Source
gf	32.21	kJ/mol	Joback Method
hf	-370.66	kJ/mol	Joback Method
hfus	38.41	kJ/mol	Joback Method
hvap	66.70	kJ/mol	Joback Method
log10ws	-4.94		Crippen Method
logp	4.694		Crippen Method
mvol	237.300	ml/mol	McGowan Method
pc	1636.45	kPa	Joback Method
rinpol	2128.00		NIST Webbook
rinpol	2128.00		NIST Webbook
tb	716.28	K	Joback Method
tc	918.20	K	Joback Method
tf	404.90	K	Joback Method
vc	0.915	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	698.47	J/molxK	716.28	Joback Method
cpg	716.99	J/molxK	749.93	Joback Method
cpg	734.39	J/molxK	783.59	Joback Method
cpg	750.74	J/molxK	817.24	Joback Method
cpg	766.08	J/molxK	850.89	Joback Method
cpg	780.47	J/molxK	884.54	Joback Method
cpg	793.96	J/molxK	918.20	Joback Method
dvisc	0.0018915	Paxs	404.90	Joback Method

dvisc	0.0011927	Paxs	456.80	Joback Method
dvisc	0.0008263	Paxs	508.69	Joback Method
dvisc	0.0006127	Paxs	560.59	Joback Method
dvisc	0.0004779	Paxs	612.49	Joback Method
dvisc	0.0003875	Paxs	664.38	Joback Method
dvisc	0.0003240	Paxs	716.28	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U405999&Units=SI
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

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
log_p:	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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