

trans-2,3-epoxydecanal

Inchi:	InChI=1S/C10H18O2/c1-2-3-4-5-6-7-9-10(8-11)12-9/h8-10H,2-7H2,1H3/t9-,10-/m0/s1
InchiKey:	ZAYNCWQQQVRPDV-UWVGGRQHSA-N
Formula:	C10H18O2
SMILES:	CCCCCCCC1OC1C=O
Mol. weight [g/mol]:	170.25

Physical Properties

Property code	Value	Unit	Source
gf	-99.28	kJ/mol	Joback Method
hf	-414.85	kJ/mol	Joback Method
hfus	31.13	kJ/mol	Joback Method
hvap	48.69	kJ/mol	Joback Method
log10ws	-2.49		Crippen Method
logp	2.313		Crippen Method
mcvol	148.340	ml/mol	McGowan Method
pc	2460.47	kPa	Joback Method
rinpol	1296.00		NIST Webbook
rinpol	1370.00		NIST Webbook
ripol	1756.00		NIST Webbook
ripol	1756.00		NIST Webbook
tb	505.88	K	Joback Method
tc	686.62	K	Joback Method
tf	284.73	K	Joback Method
vc	0.590	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	361.06	J/molxK	505.88	Joback Method
cpg	428.70	J/molxK	656.49	Joback Method
cpg	416.49	J/molxK	626.37	Joback Method
cpg	403.65	J/molxK	596.25	Joback Method
cpg	390.15	J/molxK	566.13	Joback Method
cpg	375.96	J/molxK	536.00	Joback Method

cpg	440.30	J/molxK	686.62	Joback Method
dvisc	0.0006446	Paxs	505.88	Joback Method
dvisc	0.0007423	Paxs	469.02	Joback Method
dvisc	0.0008756	Paxs	432.16	Joback Method
dvisc	0.0010650	Paxs	395.31	Joback Method
dvisc	0.0013488	Paxs	358.45	Joback Method
dvisc	0.0018032	Paxs	321.59	Joback Method
dvisc	0.0025988	Paxs	284.73	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=R237060&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
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

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
ripol:	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/76-533-6/trans-2-3-epoxydecanal.pdf>

Generated by Cheméo on 2024-04-20 14:11:09.392496858 +0000 UTC m=+15911518.313074173.
Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.