Dicyclopentadiene (endo)

Inchi: InChI=1S/C10H12/c1-2-9-7-4-5-8(6-7)10(9)3-1/h1-2,4-5,7-10H,3,6H2/t7?,8?,9-,10?/m0/s
InchiKey: HECLRDQVFMWTQS-HTEYAMCKSA-N
Formula: C10H12
SMILES: C1=CC2C3C=CC(C3)C2C1
Mol. weight [g/mol]: 132.20

Physical Properties

Property code	Value	Unit	Source
gf	255.68	kJ/mol	Joback Method
hf	57.73	kJ/mol	Joback Method
hfus	17.48	kJ/mol	Joback Method
hvap	38.04	kJ/mol	Joback Method
log10ws	-2.43		Crippen Method
logp	2.385		Crippen Method
mcvol	110.580	ml/mol	McGowan Method
pc	3364.54	kPa	Joback Method
rinpol	1032.40		NIST Webbook
rinpol	1027.00		NIST Webbook
rinpol	1017.00		NIST Webbook
rinpol	1025.00		NIST Webbook
rinpol	1016.00		NIST Webbook
rinpol	989.00		NIST Webbook
rinpol	1003.30		NIST Webbook
rinpol	987.00		NIST Webbook
rinpol	1025.00		NIST Webbook
rinpol	1005.00		NIST Webbook
rinpol	1016.00		NIST Webbook
rinpol	1043.20		NIST Webbook
rinpol	1011.90		NIST Webbook
ripol	1245.10		NIST Webbook
ripol	1247.10		NIST Webbook
ripol	1245.10		NIST Webbook
tb	446.34	K	Joback Method
tc	664.47	K	Joback Method
tf	250.04	K	Joback Method
vc	0.429	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	242.89	J/mol×K	446.34	Joback Method
cpg	261.44	J/mol×K	482.70	Joback Method
cpg	278.56	J/mol×K	519.05	Joback Method
cpg	294.37	J/mol×K	555.41	Joback Method
cpg	308.96	J/mol×K	591.76	Joback Method
cpg	322.43	J/mol×K	628.12	Joback Method
cpg	334.89	J/mol×K	664.47	Joback Method
dvisc	0.0004087	Pa×s	250.04	Joback Method
dvisc	0.0005202	Pa×s	282.76	Joback Method
dvisc	0.0006298	Pa×s	315.47	Joback Method
dvisc	0.0007356	Pa×s	348.19	Joback Method
dvisc	0.0008365	Pa×s	380.91	Joback Method
dvisc	0.0009322	Pa×s	413.62	Joback Method
dvisc	0.0010224	Pa×s	446.34	Joback Method

Sources

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
NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=R211959&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
ripol:	Polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume