ISOLONGIFOLENE, 9,10-DEHYDRO-

Other names: 9,10-Dehydroisolongifolene
Inchi: InChI=1S/C15H22/c1-13(2)8-5-6-12-14(3,4)11-7-9-15(12,13)10-11/h5-6,8,11H,7,9-10H2,
InchiKey: UIPKEVNEOFKIRG-UHFFFAOYSA-N
Formula: C15H22
SMILES: CC1(C)C2=CC=CC(C)(C)C23CCC1C3
Mol. weight [g/mol]: 202.34

Physical Properties

Property code	Value	Unit	Source
hfus	9.04	kJ/mol	Joback Method
logp	4.335		Crippen Method
mcvol	181.030	ml/mol	McGowan Method
ripol	1913.00		NIST Webbook
ripol	1913.00		NIST Webbook

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	481.77	J/mol×K	570.71	Joback Method
cpg	502.91	J/mol×K	610.26	Joback Method
cpg	522.45	J/mol×K	649.81	Joback Method
cpg	540.85	J/mol×K	689.36	Joback Method
cpg	558.58	J/mol×K	728.91	Joback Method
cpg	576.10	J/mol×K	768.46	Joback Method
cpg	593.88	J/mol×K	808.01	Joback Method

Sources

McGowan Method: http://link.springer.com/article/10.1007/BF02311772
Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws

Cheméo Relay (1.0):	
Cheméo Relay (1.0, beta):	
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U151671&Units=SI
Joback Method:	https://en.wikipedia.org/wiki/Joback_method

Legend

cpg:	Ideal gas heat capacity
hfus:	Enthalpy of fusion at standard conditions
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
ripol:	Polar retention indices