Cyclohexene, 4-[(1E)-1,5-dimethyl-1,4-hexadien-1-yl]-1-methyl-

Other names: 4-[(1E)-1,5-Dimethyl-1,4-hexadienyl]-1-methyl-1-cyclohexene
2,5-Heptadiene, 2-methyl-6-(4-methyl-3-cyclohexen-1-yl)-, (E)-
Cyclohexene, 4-(1,5-dimethyl-1,4-hexadienyl)-1-methyl-, (E)-
Cyclohexene, 4-[(1E)-1,5-dimethyl-1,4-hexadienyl]-1-methyl-
(E)-«alpha»-Bisabolene
trans-«alpha»-Bisabolene
«alpha»-Bisabolene

Inchi: InChI=1S/C15H24/c1-12(2)6-5-7-14(4)15-10-8-13(3)9-11-15/h6-8,15H,5,9-11H2,1-4H3

InchiKey: YHBUQBJHSRGZNF-UHFFFAOYSA-N

Formula: C15H24

SMILES: CC(C)=CCC=C(C)C1CC=C(C)CC1

Mol. weight [g/mol]: 204.35

CAS: 25532-79-0

Physical Properties

Property code	Value	Unit	Source
gf	263.54	kJ/mol	Joback Method
hf	-37.44	kJ/mol	Joback Method
hfus	25.06	kJ/mol	Joback Method
hvap	50.44	kJ/mol	Joback Method
log10ws	-5.32		Crippen Method
logp	5.035		Crippen Method
mcvol	198.450	ml/mol	McGowan Method
pc	1861.11	kPa	Joback Method
rinpol	1504.00		NIST Webbook
rinpol	1491.00		NIST Webbook
rinpol	1532.00		NIST Webbook
rinpol	1506.00		NIST Webbook
rinpol	1529.00		NIST Webbook
rinpol	1493.00		NIST Webbook
rinpol	1531.00		NIST Webbook
rinpol	1493.00		NIST Webbook
rinpol	1531.00		NIST Webbook
rinpol	1496.00		NIST Webbook
rinpol	1496.00		NIST Webbook
rinpol	1496.00		NIST Webbook
rinpol	1544.60		NIST Webbook

rinpol	1511.00	NIST Webbook
rinpol	1506.00	NIST Webbook
rinpol	1523.00	NIST Webbook
rinpol	1531.00	NIST Webbook
rinpol	1531.00	NIST Webbook
rinpol	1508.00	NIST Webbook
rinpol	1528.00	NIST Webbook
rinpol	1487.00	NIST Webbook
rinpol	1500.00	NIST Webbook
rinpol	1496.00	NIST Webbook
rinpol	1503.00	NIST Webbook
rinpol	1515.00	NIST Webbook
rinpol	1512.00	NIST Webbook
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rinpol	1500.00	NIST Webbook
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rinpol	1499.00		NIST Webbook
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ripol	1766.00		NIST Webbook
ripol	1736.00		NIST Webbook
ripol	1777.00		NIST Webbook
ripol	1737.50		NIST Webbook
ripol	1777.00		NIST Webbook
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ripol	1720.00		NIST Webbook
ripol	1724.00		NIST Webbook
ripol	1766.00		NIST Webbook
ripol	1733.00		NIST Webbook
ripol	1776.00		NIST Webbook
ripol	1720.00		NIST Webbook
ripol	1786.00		NIST Webbook
ripol	1770.00		NIST Webbook
ripol	1724.00		NIST Webbook
ripol	1726.00		NIST Webbook
tb	574.37	K	Joback Method
tc	785.97	K	Joback Method
tf	241.39	K	Joback Method
vc	0.756	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	495.01	J/mol×K	574.37	Joback Method
cpg	516.16	J/mol×K	609.64	Joback Method
cpg	536.07	J/mol×K	644.90	Joback Method
cpg	554.80	J/mol×K	680.17	Joback Method
cpg	572.42	J/mol×K	715.43	Joback Method
cpg	589.00	J/mol×K	750.70	Joback Method
cpg	604.60	J/mol×K	785.97	Joback Method

Sources

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=C25532790&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws

Legend

cpg:	Ideal gas heat capacity
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