

Indan, 2-butyl-5-hexyl-

Other names:	2-n-Butyl-5-n-hexyl-(2,3-dihydroindene) 2-n-Butyl-5-n-hexylindan
Inchi:	InChI=1S/C19H30/c1-3-5-7-8-10-16-11-12-18-14-17(9-6-4-2)15-19(18)13-16/h11-13,17H
InchiKey:	JAUZOZVMGVOYCU-UHFFFAOYSA-N
Formula:	C19H30
SMILES:	CCCCCc1ccc2c(c1)CC(CCCC)C2
Mol. weight [g/mol]:	258.44
CAS:	25446-32-6

Physical Properties

Property code	Value	Unit	Source
gf	263.00	kJ/mol	Joback Method
hf	-149.10	kJ/mol	Joback Method
hfus	36.36	kJ/mol	Joback Method
hvap	61.40	kJ/mol	Joback Method
log10ws	-6.46		Crippen Method
logp	5.714		Crippen Method
mvol	243.950	ml/mol	McGowan Method
pc	1466.85	kPa	Joback Method
tb	677.50	K	Joback Method
tc	873.70	K	Joback Method
tf	239.45 ± 0.50	K	NIST Webbook
vc	0.949	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	696.51	J/mol×K	677.50	Joback Method
cpg	787.03	J/mol×K	841.00	Joback Method
cpg	770.86	J/mol×K	808.30	Joback Method
cpg	753.78	J/mol×K	775.60	Joback Method
cpg	735.74	J/mol×K	742.90	Joback Method
cpg	716.67	J/mol×K	710.20	Joback Method
cpg	802.35	J/mol×K	873.70	Joback Method

dvisc	0.0002868	Paxs	677.50	Joback Method
dvisc	0.0003436	Paxs	626.80	Joback Method
dvisc	0.0004250	Paxs	576.10	Joback Method
dvisc	0.0005478	Paxs	525.39	Joback Method
dvisc	0.0007453	Paxs	474.69	Joback Method
dvisc	0.0010915	Paxs	423.99	Joback Method
dvisc	0.0017731	Paxs	373.29	Joback Method

Sources

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=C25446326&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071

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
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/41-787-3/Indan-2-butyl-5-hexyl.pdf>

Generated by Cheméo on 2024-04-19 17:12:43.702322511 +0000 UTC m=+15836012.622899822.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.