

2,6,10-trimethyl-7-(3-methylbutyl)dodecane

Inchi:	InChI=1S/C20H42/c1-8-18(6)13-15-20(14-12-17(4)5)19(7)11-9-10-16(2)3/h16-20H,8-15H
InchiKey:	LSVDXMJHSYNAGC-UHFFFAOYSA-N
Formula:	C20H42
SMILES:	CCC(C)CCC(CCC(C)C)C(C)CCCC(C)C
Mol. weight [g/mol]:	282.55

Physical Properties

Property code	Value	Unit	Source
gf	105.32	kJ/mol	Joback Method
hf	-482.53	kJ/mol	Joback Method
hfus	29.94	kJ/mol	Joback Method
hvap	58.17	kJ/mol	Joback Method
log10ws	-6.99		Crippen Method
logp	7.328		Crippen Method
mcvol	292.660	ml/mol	McGowan Method
pc	1041.93	kPa	Joback Method
rinpol	1706.00		NIST Webbook
rinpol	1706.00		NIST Webbook
tb	654.80	K	Joback Method
tc	824.55	K	Joback Method
tf	240.16	K	Joback Method
vc	1.125	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	842.14	J/mol×K	654.80	Joback Method
cpg	864.47	J/mol×K	683.09	Joback Method
cpg	885.81	J/mol×K	711.38	Joback Method
cpg	906.19	J/mol×K	739.68	Joback Method
cpg	925.64	J/mol×K	767.97	Joback Method
cpg	944.18	J/mol×K	796.26	Joback Method
cpg	961.85	J/mol×K	824.55	Joback Method
dvisc	0.0359131	Paxs	240.16	Joback Method

dvisc	0.0039123	Paxs	309.27	Joback Method
dvisc	0.0009579	Paxs	378.37	Joback Method
dvisc	0.0003622	Paxs	447.48	Joback Method
dvisc	0.0001777	Paxs	516.59	Joback Method
dvisc	0.0001031	Paxs	585.69	Joback Method
dvisc	0.0000671	Paxs	654.80	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=R261939&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
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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