

# Benzene, 1-hexyl-2-propyl

<b>Inchi:</b>	InChI=1S/C15H24/c1-3-5-6-7-11-15-13-9-8-12-14(15)10-4-2/h8-9,12-13H,3-7,10-11H2,1
<b>InchiKey:</b>	RVFXOHWJVXHUHE-UHFFFAOYSA-N
<b>Formula:</b>	C15H24
<b>SMILES:</b>	CCCCCc1ccccc1CCC
<b>Mol. weight [g/mol]:</b>	204.35

## Physical Properties

Property code	Value	Unit	Source
gf	178.20	kJ/mol	Joback Method
hf	-127.87	kJ/mol	Joback Method
hfus	28.26	kJ/mol	Joback Method
hvap	51.92	kJ/mol	Joback Method
log10ws	-5.17		Crippen Method
logp	4.762		Crippen Method
mvol	198.450	ml/mol	McGowan Method
pc	1823.17	kPa	Joback Method
tb	574.26	K	Joback Method
tc	766.81	K	Joback Method
tf	297.75	K	Joback Method
vc	0.767	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	493.06	J/molxK	574.26	Joback Method
cpg	511.45	J/molxK	606.35	Joback Method
cpg	528.91	J/molxK	638.44	Joback Method
cpg	545.48	J/molxK	670.54	Joback Method
cpg	561.18	J/molxK	702.63	Joback Method
cpg	576.06	J/molxK	734.72	Joback Method
cpg	590.13	J/molxK	766.81	Joback Method
dvisc	0.0025521	Paxs	297.75	Joback Method
dvisc	0.0011846	Paxs	343.83	Joback Method
dvisc	0.0006592	Paxs	389.92	Joback Method

dvisc	0.0004153	Paxs	436.00	Joback Method
dvisc	0.0002857	Paxs	482.09	Joback Method
dvisc	0.0002099	Paxs	528.17	Joback Method
dvisc	0.0001620	Paxs	574.26	Joback Method

## Sources

<b>Joback Method:</b>	<a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>
<b>McGowan Method:</b>	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>
<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=R13692&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R13692&amp;Units=SI</a>
<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>
<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>

## Legend

<b>cpg:</b>	Ideal gas heat capacity
<b>dvisc:</b>	Dynamic viscosity
<b>gf:</b>	Standard Gibbs free energy of formation
<b>hf:</b>	Enthalpy of formation at standard conditions
<b>hfus:</b>	Enthalpy of fusion at standard conditions
<b>hvap:</b>	Enthalpy of vaporization at standard conditions
<b>log10ws:</b>	Log10 of Water solubility in mol/l
<b>logp:</b>	Octanol/Water partition coefficient
<b>mcvol:</b>	McGowan's characteristic volume
<b>pc:</b>	Critical Pressure
<b>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature
<b>tf:</b>	Normal melting (fusion) point
<b>vc:</b>	Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/75-092-7/Benzene-1-hexyl-2-propyl.pdf>

Generated by Cheméo on 2024-04-26 14:46:43.177357886 +0000 UTC m=+16432052.097935201.

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