

curcumenyl alcohol

Other names:	ar-Curcumen-15-ol
Inchi:	InChI=1S/C15H22O/c1-12(2)5-4-6-13(3)15-9-7-14(11-16)8-10-15/h5,7-10,13,16H,4,6,11
InchiKey:	FVVUIXKKLJOYDU-UHFFFAOYSA-N
Formula:	C15H22O
SMILES:	CC(C)=CCCC(C)c1ccc(CO)cc1
Mol. weight [g/mol]:	218.33

Physical Properties

Property code	Value	Unit	Source
gf	110.61	kJ/mol	Joback Method
hf	-177.95	kJ/mol	Joback Method
hfus	27.71	kJ/mol	Joback Method
hvap	68.25	kJ/mol	Joback Method
log10ws	-4.75		Crippen Method
logp	4.029		Crippen Method
mcvol	200.020	ml/mol	McGowan Method
pc	2094.58	kPa	Joback Method
rinpol	1705.00		NIST Webbook
tb	670.04	K	Joback Method
tc	865.73	K	Joback Method
tf	324.53	K	Joback Method
vc	0.761	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	538.91	J/mol×K	670.04	Joback Method
cpg	554.16	J/mol×K	702.66	Joback Method
cpg	568.57	J/mol×K	735.27	Joback Method
cpg	582.17	J/mol×K	767.89	Joback Method
cpg	595.02	J/mol×K	800.50	Joback Method
cpg	607.16	J/mol×K	833.12	Joback Method
cpg	618.65	J/mol×K	865.73	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=R314028&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l

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
rinpola:	Non-polar retention indices
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/79-027-5/curcumenyl-alcohol.pdf>

Generated by Cheméo on 2025-12-06 01:23:57.252963983 +0000 UTC m=+4732434.783004638.

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