

# 4-Heptylbenzoic acid

<b>Other names:</b>	Benzoic acid, 4-heptyl-p-heptylbenzoic acid
<b>Inchi:</b>	InChI=1S/C14H20O2/c1-2-3-4-5-6-7-12-8-10-13(11-9-12)14(15)16/h8-11H,2-7H2,1H3,(H
<b>InchiKey:</b>	VSUKEWPHURLYTK-UHFFFAOYSA-N
<b>Formula:</b>	C14H20O2
<b>SMILES:</b>	CCCCCCCc1ccc(C(=O)O)cc1
<b>Mol. weight [g/mol]:</b>	220.31
<b>CAS:</b>	38350-87-7

## Physical Properties

Property code	Value	Unit	Source
gf	-95.96	kJ/mol	Joback Method
hf	-372.04	kJ/mol	Joback Method
hfus	1.00	kJ/mol	Thermodynamic study of the sublimation of eight 4-n-alkylbenzoic acids
hsub	130.00 ± 0.90	kJ/mol	NIST Webbook
hvap	73.12	kJ/mol	Joback Method
log10ws	-4.43		Crippen Method
logp	3.898		Crippen Method
mcvol	191.800	ml/mol	McGowan Method
pc	2287.14	kPa	Joback Method
tb	697.43	K	Joback Method
tc	889.73	K	Joback Method
tf	376.00 ± 1.00	K	NIST Webbook
vc	0.737	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	527.16	J/mol×K	697.43	Joback Method
cpg	540.70	J/mol×K	729.48	Joback Method
cpg	553.48	J/mol×K	761.53	Joback Method
cpg	565.53	J/mol×K	793.58	Joback Method

cpg	576.88	J/molxK	825.63	Joback Method
cpg	587.57	J/molxK	857.68	Joback Method
cpg	597.61	J/molxK	889.73	Joback Method
dvisc	0.0022271	Paxs	397.23	Joback Method
dvisc	0.0008134	Paxs	447.26	Joback Method
dvisc	0.0003638	Paxs	497.30	Joback Method
dvisc	0.0001885	Paxs	547.33	Joback Method
dvisc	0.0001091	Paxs	597.36	Joback Method
dvisc	0.0000687	Paxs	647.40	Joback Method
dvisc	0.0000462	Paxs	697.43	Joback Method

## Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.20363e+01
Coeff. B	-3.91173e+03
Coeff. C	-9.40500e+01
Temperature range (K), min.	427.00
Temperature range (K), max.	675.73

## Sources

Crippen Method:

[https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)

Thermodynamic study of the sublimation of eight 4-n-alkylbenzoic acids

<https://www.doi.org/10.1016/j.jct.2004.02.001>

Joback Method:

[https://en.wikipedia.org/wiki/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=C38350877&Units=SI>

The Yaws Handbook of Vapor Pressure:  
Crippen Method:

<https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure>

<http://pubs.acs.org/doi/abs/10.1021/ci9903071>

## Legend

cpg: Ideal gas heat capacity

dvisc: 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>hsub:</b>	Enthalpy of sublimation 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>pvap:</b>	Vapor 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

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