

# Benzenethiol, 2-methyl-

<b>Other names:</b>	2-Methylbenzenethiol 2-Methylbenzothiol 2-Methylphenylthiol 2-Methylthiophenol 2-Toluenethiol NSC 80655 USAF EK-2676 o-Mercaptotoluene o-Methylbenzenethiol o-Methylthiophenol o-Thiocresol o-Toluenethiol o-Tolyl mercaptan toluene-2-thiol
<b>Inchi:</b>	InChI=1S/C7H8S/c1-6-4-2-3-5-7(6)8/h2-5,8H,1H3
<b>InchiKey:</b>	LXUNZSDDXMPKLP-UHFFFAOYSA-N
<b>Formula:</b>	C7H8S
<b>SMILES:</b>	Cc1ccccc1S
<b>Mol. weight [g/mol]:</b>	124.20
<b>CAS:</b>	137-06-4

## Physical Properties

Property code	Value	Unit	Source
gf	140.23	kJ/mol	Joback Method
hf	75.73	kJ/mol	Joback Method
hfus	11.58	kJ/mol	Joback Method
hvap	40.85	kJ/mol	Joback Method
ie	8.31	eV	NIST Webbook
log10ws	-2.46		Crippen Method
logp	2.284		Crippen Method
mcvol	102.080	ml/mol	McGowan Method
pc	4385.77	kPa	Joback Method
tb	468.20	K	NIST Webbook
tc	694.66	K	Joback Method
tf	244.05	K	Joback Method
vc	0.373	m3/kmol	Joback Method

# Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	181.55	J/mol×K	454.08	Joback Method
cpg	192.97	J/mol×K	494.18	Joback Method
cpg	203.66	J/mol×K	534.27	Joback Method
cpg	213.64	J/mol×K	574.37	Joback Method
cpg	222.96	J/mol×K	614.47	Joback Method
cpg	231.63	J/mol×K	654.56	Joback Method
cpg	239.70	J/mol×K	694.66	Joback Method
hvapt	48.10	kJ/mol	424.50	NIST Webbook
hvapt	46.60	kJ/mol	420.00	NIST Webbook

## Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.45809e+01
Coeff. B	-4.10903e+03
Coeff. C	-5.57550e+01
Temperature range (K), min.	343.23
Temperature range (K), max.	499.04

## Sources

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>

Crippen Method:

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

Joback Method:

[https://en.wikipedia.org/wiki/Joback\\_method](https://en.wikipedia.org/wiki/Joback_method)

KDB:

<https://www.thermo.com/files/research/kdb/mol/mol1854.mol>

McGowan Method:

<http://link.springer.com/article/10.1007/BF02311772>

NIST Webbook:

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C137064&Units=SI>

# Legend

<b>cpg:</b>	Ideal gas heat capacity
<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>hvapt:</b>	Enthalpy of vaporization at a given temperature
<b>ie:</b>	Ionization energy
<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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