

# Allylphenyl sulfide

<b>Other names:</b>	(Allylthio)benzene Benzene, (2-propenylthio)- 3-(Phenylthio)propene 3-(Phenylthio)prop-1-ene Sulfide, allyl phenyl Allyl phenyl thioether Phenyl allyl thioether Phenyl allyl sulfide
<b>Inchi:</b>	InChI=1S/C9H10S/c1-2-8-10-9-6-4-3-5-7-9/h2-7H,1,8H2
<b>InchiKey:</b>	QGNRLAFFKKBSIM-UHFFFAOYSA-N
<b>Formula:</b>	C9H10S
<b>SMILES:</b>	C=CCSc1ccccc1
<b>Mol. weight [g/mol]:</b>	150.24
<b>CAS:</b>	5296-64-0

## Physical Properties

Property code	Value	Unit	Source
gf	258.27	kJ/mol	Joback Method
hf	174.74	kJ/mol	Joback Method
hfus	15.96	kJ/mol	Joback Method
hvap	44.05	kJ/mol	Joback Method
ie	7.91 ± 0.01	eV	NIST Webbook
log10ws	-2.91		Crippen Method
logp	2.965		Crippen Method
mcvol	125.960	ml/mol	McGowan Method
pc	3493.01	kPa	Joback Method
tb	497.46	K	Joback Method
tc	733.42	K	Joback Method
tf	250.25	K	Joback Method
vc	0.467	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
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cpg	247.62	J/mol×K	497.46	Joback Method
cpg	261.28	J/mol×K	536.79	Joback Method
cpg	274.02	J/mol×K	576.11	Joback Method
cpg	285.86	J/mol×K	615.44	Joback Method
cpg	296.85	J/mol×K	654.77	Joback Method
cpg	307.04	J/mol×K	694.09	Joback Method
cpg	316.46	J/mol×K	733.42	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=C5296640&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C5296640&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>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>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>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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