

# S-diphenylcarbazone

**Inchi:** InChI=1S/C26H24N8O2/c35-25(31-27-21-7-3-1-4-8-21)33-29-23-15-11-19(12-16-23)20-  
**InchiKey:** LEKPXSOMGUDVOS-VBMGMRCRSA-N  
**Formula:** C26H24N8O2  
**SMILES:** O=C(N=Nc1ccc(-c2ccc(NNC(=O)NNc3ccccc3)cc2)cc1)NNc1ccccc1  
**Mol. weight [g/mol]:** 480.52

## Physical Properties

Property code	Value	Unit	Source
hf	486.09	kJ/mol	Joback Method
hvap	142.68	kJ/mol	Joback Method
log10ws	-8.85		Crippen Method
logp	5.827		Crippen Method
mcvol	360.840	ml/mol	McGowan Method
pc	1628.54	kPa	Joback Method
tb	1468.92	K	Joback Method
tc	1805.03	K	Joback Method

## Sources

**McGowan Method:** <http://link.springer.com/article/10.1007/BF02311772>  
**NIST Webbook:** <http://webbook.nist.gov/cgi/cbook.cgi?ID=B6000377&Units=SI>  
**Crippen Method:** <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)

## Legend

**hf:** Enthalpy of formation 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  
**tb:** Normal Boiling Point Temperature  
**tc:** Critical Temperature

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