

# Thiabendazole

**Other names:** 1H-Benzimidazole, 2-(4-thiazolyl)-  
Benzimidazole, 2-(4-thiazolyl)-  
Equizole  
Mertec  
Mertect  
Mertect 160  
Metasol TK 100  
Mintesol  
Mintezol  
Minzolum  
MK 360  
Omnizole  
Tebuzate  
Tecto  
Tecto 10P  
Tecto 40F  
Tecto 60  
Testo  
Thiabendazol  
Thiabendazole  
Thibenzol  
Thibenzole  
Thibenzole 200  
Tiabenda  
Triasox  
TBZ  
TBZ 6  
TBZ 60W  
2-(Thiazol-4-yl)benzimidazole  
2-(4-Thiazolyl)benzimidazole  
2-(4'-Thiazolyl)benzimidazole  
Apl-Luster  
Bovizole  
E-Z-Ex  
Eprofil  
Lombristop  
Mycozol  
Nemapan  
Polival  
Thiaben

Thibenzole ATT  
Top form wormer  
2-Thiazole-4-ylbenzimidazole  
Arbotect  
4-(2-Benzimidazolyl)thiazole  
Bioguard  
RPH  
Storite  
Tbdz  
Tecto rph  
Thiabenzazole  
2-(4-Thiazolyl)-1H-benzimidazole  
Tobaz  
Tiabendazole  
Equivet TZ  
Tiabendazol  
Tibimix 20  
Thibendole  
Thiabendole  
Sistemas  
Sanaizol 100  
Chemviron TK 100  
Tecto B  
5-(4-Thiazolyl)benzimidazole  
Ormogal  
Cropasal  
Drawipas  
g 491  
Hokustar HP  
Mertect 340F  
Mertect LSP  
2-(1,3-Thiazol-4-yl)benzimidazole  
Hymush  
Nemacin  
Thiprazole  
Tubazole  
2-(4-Thiazolyl)-1H-benzimidazole  
4-(1H-Benzimidazol-2-yl)thiazole  
Pitrizet  
Metasol TK 25  
Statin  
2-(4-Thiazolyl)-1H-benzimidazole (tiabendazole)  
**Inchi:** InChI=1S/C10H7N3S/c1-2-4-8-7(3-1)12-10(13-8)9-5-14-6-11-9/h1-6H,(H,12,13)

**InchiKey:** WJCNZQLZVWNLKY-UHFFFAOYSA-N  
**Formula:** C10H7N3S  
**SMILES:** c1ccc2[nH]c(-c3cscn3)nc2c1  
**Mol. weight [g/mol]:** 201.25  
**CAS:** 148-79-8

## Physical Properties

Property code	Value	Unit	Source
log10ws	-4.39		Crippen Method
logp	2.204		Crippen Method
mcvol	139.670	ml/mol	McGowan Method
rinpol	2088.00		NIST Webbook
rinpol	2080.00		NIST Webbook
rinpol	2010.00		NIST Webbook
rinpol	2091.00		NIST Webbook
rinpol	2040.00		NIST Webbook
rinpol	2080.00		NIST Webbook
rinpol	2010.00		NIST Webbook
rinpol	2091.00		NIST Webbook
tf	574.80 ± 0.20	K	NIST Webbook

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
hfust	35.20	kJ/mol	573.20	NIST Webbook

## Sources

**Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci990307l>  
**Crippen Method:** [https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)  
**McGowan Method:** <http://link.springer.com/article/10.1007/BF02311772>  
**NIST Webbook:** <http://webbook.nist.gov/cgi/cbook.cgi?ID=C148798&Units=SI>

# Legend

<b>hfust:</b>	Enthalpy of fusion at a given temperature
<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>rinpola:</b>	Non-polar retention indices
<b>tf:</b>	Normal melting (fusion) point

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