

2,2,5-endo,6-exo,8,9,9,10-octachlorobornane

Other names:	2,2,5,6-Tetrachloro-1,7-bis(chloromethyl)-7-(dichloromethyl)bicyclo[2.2.1]heptane, (5-endo,6-exo,7-syn)-2,2,5-endo,6-exo,8,8,9,10-octachlorobornane 2,2,5-endo,6-exo,8c,9b,9c,10a-octachlorobornane Bicyclo[2.2.1]heptane, 2,2,5,6-tetrachloro-1,7-bis(chloromethyl)-7-(dichloromethyl)-, (5-endo,6-exo,7-syn)-rel-(1R,4R,5R,6R,7R)-2,2,5,6-tetrachloro-1,7-bis(chloromethyl)-7-(dichloromethyl)bicyclo[2.2.1]heptane, 420
Inchi:	InChI=1S/C10H10Cl8/c11-2-8(7(15)16)4-1-10(17,18)9(8,3-12)6(14)5(4)13/h4-7H,1-3H2/t
InchiKey:	YNEKMCSWRMRXIR-UHFFFAOYSA-N
Formula:	C10H10Cl8
SMILES:	C1CC1(C(Cl)Cl)C2CC(Cl)(Cl)C1(CCl)C(Cl)C2Cl
Mol. weight [g/mol]:	413.81

Physical Properties

Property code	Value	Unit	Source
hfus	31.27	kJ/mol	Joback Method
logp	5.663		Crippen Method
mcvol	227.960	ml/mol	McGowan Method
rinpol	2373.20		NIST Webbook
rinpol	2373.20		NIST Webbook

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	508.07	J/mol×K	726.99	Joback Method
cpg	520.99	J/mol×K	770.56	Joback Method
cpg	534.50	J/mol×K	814.13	Joback Method
cpg	549.15	J/mol×K	857.69	Joback Method
cpg	565.50	J/mol×K	901.26	Joback Method
cpg	584.12	J/mol×K	944.83	Joback Method
cpg	605.56	J/mol×K	988.40	Joback Method
pvap	1.51e-06	kPa	298.15	Vapor Pressures and Enthalpies of Vaporization for Toxaphene Congeners

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Cheméo Relay (1.0):	
Cheméo Relay (1.0, beta):	
Vapor Pressures and Enthalpies of Vaporization for Toxaphene Congeners	https://www.doi.org/10.1021/je020204q
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C58002190&Units=SI
Joback Method:	https://en.wikipedia.org/wiki/Joback_method
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772

Legend

cpg:	Ideal gas heat capacity
hfus:	Enthalpy of fusion at standard conditions
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pvap:	Vapor pressure
rinpol:	Non-polar retention indices

Latest version available from:

<https://www.chemeo.com/cid/85-982-8/2-2-5-endo-6-exo-8-9-9-10-octachlorobornane.pdf>

Generated by Cheméo on 2026-07-22 02:49:48.311805856 +0000 UTC m=+199684.153691232.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.