

Data Sheet

NAME:	Indinavir Sulphate
REPOSITORY REFERENCE:	ARP972
PROVIDED:	20 mg
CHEMICAL NAME:	(alphaR, gammas,2S)-alpha-Benzyl-2-(tert-butylcarbamoyl)-gamma-hydroxy-N-[(1S,2R)-2-hydroxy-1-indanyl]-4-(3-pyridyl-methyl)-1-piperazinevaleramide sulfate.
MOLECULAR FORMULA:	$C_{36}H_{47}N_5O_4 \bullet H_2SO_4$
MOLECULAR WEIGHT:	711.88
CHARACTERISTICS:	Indinavir sulphate is a selective, competitive inhibitor of HIV-1 protease. Indinavir is a structural analogue of the HIV <i>Phe-Pro</i> protease cleavage site. The drug's structure inhibits the function of HIV protease, blocking virus maturation and causing the formation of immature, noninfectious virions.
SOLUBILITY:	Very soluble in water and in methanol. Insoluble in hexanes
RECOMMENDED STORAGE:	Room temperature

CONTRIBUTOR:

Division of AIDS, NIAID.

ACKNOWLEDGEMENT:

Publications should acknowledge the donor of the reagent and the Programme EVA Centre for AIDS Reagents. Suggested wording can be found on our website at <http://www.nibsc.ac.uk/spotlight/aidsreagent/index.html> in the "Acknowledgements" section.

Please also ensure that you send us a copy of any papers resulting from work using reagents acquired through CFAR (this can be electronically or as a paper copy)

NOTE:

This compound is limited to 40 mg per requester per year. **Recipient agrees that the reagent (Indinavir Sulphate) use is permitted only as a standard for in vitro and/or studies in animals for inhibition of HIV replication.**