



# wwPDB X-ray Structure Validation Summary Report ⓘ

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PDB ID : 1NB8  
Title : Structure of the catalytic domain of USP7 (HAUSP)  
Authors : Hu, M.; Li, P.; Li, M.; Li, W.; Yao, T.; Wu, J.-W.; Gu, W.; Cohen, R.E.; Shi, Y.  
Deposited on : 2002-12-02  
Resolution : 2.30 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

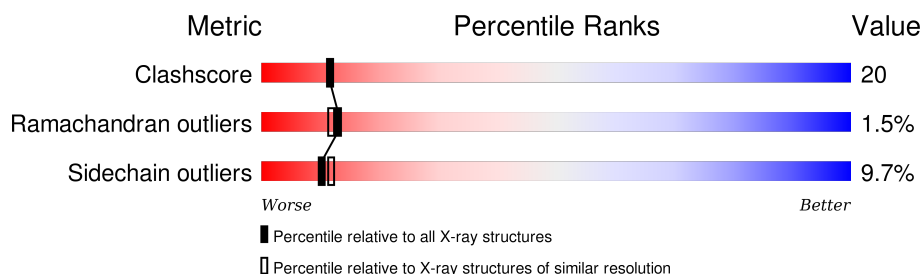
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	353	
1	B	353	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5714 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Ubiquitin carboxyl-terminal hydrolase 7.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	332	Total	C	N	O	S	Se	0	0	0
			2696	1709	461	510	7	9			
1	B	333	Total	C	N	O	S	Se	0	0	0
			2707	1715	465	511	7	9			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	225	MSE	MET	MODIFIED RESIDUE	UNP Q93009
A	244	MSE	MET	MODIFIED RESIDUE	UNP Q93009
A	245	MSE	MET	MODIFIED RESIDUE	UNP Q93009
A	292	MSE	MET	MODIFIED RESIDUE	UNP Q93009
A	311	MSE	MET	MODIFIED RESIDUE	UNP Q93009
A	328	MSE	MET	MODIFIED RESIDUE	UNP Q93009
A	407	MSE	MET	MODIFIED RESIDUE	UNP Q93009
A	410	MSE	MET	MODIFIED RESIDUE	UNP Q93009
A	515	MSE	MET	MODIFIED RESIDUE	UNP Q93009
B	225	MSE	MET	MODIFIED RESIDUE	UNP Q93009
B	244	MSE	MET	MODIFIED RESIDUE	UNP Q93009
B	245	MSE	MET	MODIFIED RESIDUE	UNP Q93009
B	292	MSE	MET	MODIFIED RESIDUE	UNP Q93009
B	311	MSE	MET	MODIFIED RESIDUE	UNP Q93009
B	328	MSE	MET	MODIFIED RESIDUE	UNP Q93009
B	407	MSE	MET	MODIFIED RESIDUE	UNP Q93009
B	410	MSE	MET	MODIFIED RESIDUE	UNP Q93009
B	515	MSE	MET	MODIFIED RESIDUE	UNP Q93009

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	139	Total	O	0	0
			139	139		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	172	Total	O	0	0
			172	172		



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	75.65Å 68.51Å 76.26Å 90.00° 95.36° 90.00°	Depositor
Resolution (Å)	20.00 – 2.30	Depositor
% Data completeness (in resolution range)	(Not available) (20.00-2.30)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	CNS	Depositor
R, $R_{free}$	0.222 , 0.279	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5714	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	47.0	wwPDB-VP

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	0.42	0/2743	0.71	4/3679 (0.1%)
1	B	0.45	1/2754 (0.0%)	0.72	0/3693
All	All	0.43	1/5497 (0.0%)	0.71	4/7372 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	223	CYS	CB-SG	5.04	1.90	1.82

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	210	HIS	N-CA-C	6.28	127.96	111.00
1	A	510	CYS	N-CA-C	6.07	127.39	111.00
1	A	235	THR	N-CA-C	-5.60	95.88	111.00
1	A	463	GLY	N-CA-C	5.15	125.97	113.10

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2696	0	2649	136	0
1	B	2707	0	2662	81	0
2	A	139	0	0	14	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	172	0	0	13	0
All	All	5714	0	5311	217	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 20.

The worst 5 of 217 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:244:MSE:HG2	2:B:563:HOH:O	1.62	0.99
1:A:419:ILE:HG22	1:A:420:LYS:H	1.32	0.95
1:B:314:THR:HG22	1:B:316:VAL:H	1.32	0.94
1:B:410:MSE:HE2	1:B:421:ILE:HD11	1.49	0.93
1:A:359:ASN:ND2	1:A:361:PHE:HB3	1.86	0.91

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	326/353 (92%)	293 (90%)	26 (8%)	7 (2%)	9	7
1	B	327/353 (93%)	304 (93%)	20 (6%)	3 (1%)	21	24
All	All	653/706 (92%)	597 (91%)	46 (7%)	10 (2%)	13	12

5 of 10 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	460	ASN
1	A	510	CYS
1	B	460	ASN

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Mol	Chain	Res	Type
1	A	375	GLY
1	A	407	MSE

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	299 / 311 (96%)	270 (90%)	29 (10%)	10	12
1	B	300 / 311 (96%)	271 (90%)	29 (10%)	10	12
All	All	599 / 622 (96%)	541 (90%)	58 (10%)	10	12

5 of 58 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	496	HIS
1	B	238	LEU
1	B	496	HIS
1	A	524	LEU
1	B	211	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 29 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	534	HIS
1	B	210	HIS
1	B	470	ASN
1	A	539	GLN
1	B	230	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

### 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers [i](#)

EDS was not executed - this section will therefore be empty.