



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 14, 2017 – 06:51 pm GMT

PDB ID : 3VLF
Title : Crystal structure of yeast proteasome interacting protein
Authors : Takagi, K.; Kim, S.; Kato, K.; Tanaka, K.; Saeki, Y.; Mizushima, T.
Deposited on : 2011-12-01
Resolution : 3.80 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

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	recalc28949

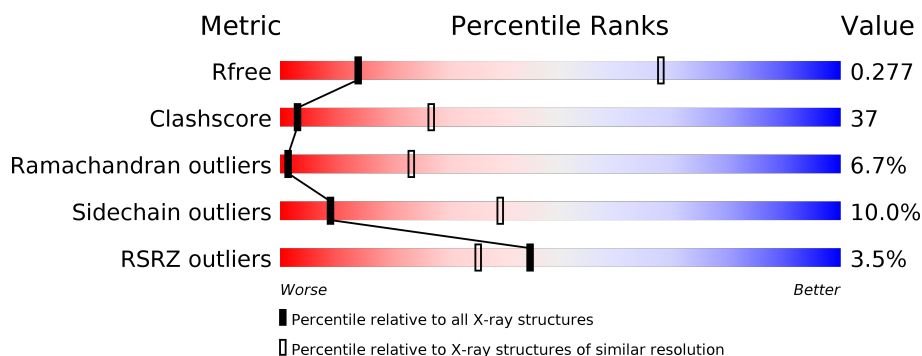
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.80 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))
R_{free}	100719	1019 (4.08-3.52)
Clashscore	112137	1030 (4.04-3.56)
Ramachandran outliers	110173	1011 (4.06-3.54)
Sidechain outliers	110143	1005 (4.06-3.54)
RSRZ outliers	101464	1032 (4.08-3.52)

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 ≥ 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.

Mol	Chain	Length	Quality of chain
1	A	500	<div> <div>5%</div> <div> <div>38%</div> <div>47%</div> <div>7%</div> <div>9%</div> </div> </div>
1	C	500	<div> <div>2%</div> <div> <div>31%</div> <div>49%</div> <div>9%</div> <div>9%</div> </div> </div>
2	B	88	<div> <div>32%</div> <div>41%</div> <div>10%</div> <div>17%</div> </div>
2	D	88	<div> <div>%</div> <div> <div>23%</div> <div>49%</div> <div>9%</div> <div>17%</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8594 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 DNA mismatch repair protein HSM3.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	457	Total	C	N	O	S	Se	0	0	0
			3727	2403	596	715	5	8			
1	C	454	Total	C	N	O	S	Se	0	0	0
			3717	2396	594	714	5	8			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MSE	-	EXPRESSION TAG	UNP P38348
A	-18	GLY	-	EXPRESSION TAG	UNP P38348
A	-17	SER	-	EXPRESSION TAG	UNP P38348
A	-16	SER	-	EXPRESSION TAG	UNP P38348
A	-15	HIS	-	EXPRESSION TAG	UNP P38348
A	-14	HIS	-	EXPRESSION TAG	UNP P38348
A	-13	HIS	-	EXPRESSION TAG	UNP P38348
A	-12	HIS	-	EXPRESSION TAG	UNP P38348
A	-11	HIS	-	EXPRESSION TAG	UNP P38348
A	-10	HIS	-	EXPRESSION TAG	UNP P38348
A	-9	SER	-	EXPRESSION TAG	UNP P38348
A	-8	SER	-	EXPRESSION TAG	UNP P38348
A	-7	GLY	-	EXPRESSION TAG	UNP P38348
A	-6	LEU	-	EXPRESSION TAG	UNP P38348
A	-5	VAL	-	EXPRESSION TAG	UNP P38348
A	-4	PRO	-	EXPRESSION TAG	UNP P38348
A	-3	ARG	-	EXPRESSION TAG	UNP P38348
A	-2	GLY	-	EXPRESSION TAG	UNP P38348
A	-1	SER	-	EXPRESSION TAG	UNP P38348
A	0	HIS	-	EXPRESSION TAG	UNP P38348
C	-19	MSE	-	EXPRESSION TAG	UNP P38348
C	-18	GLY	-	EXPRESSION TAG	UNP P38348
C	-17	SER	-	EXPRESSION TAG	UNP P38348
C	-16	SER	-	EXPRESSION TAG	UNP P38348
C	-15	HIS	-	EXPRESSION TAG	UNP P38348

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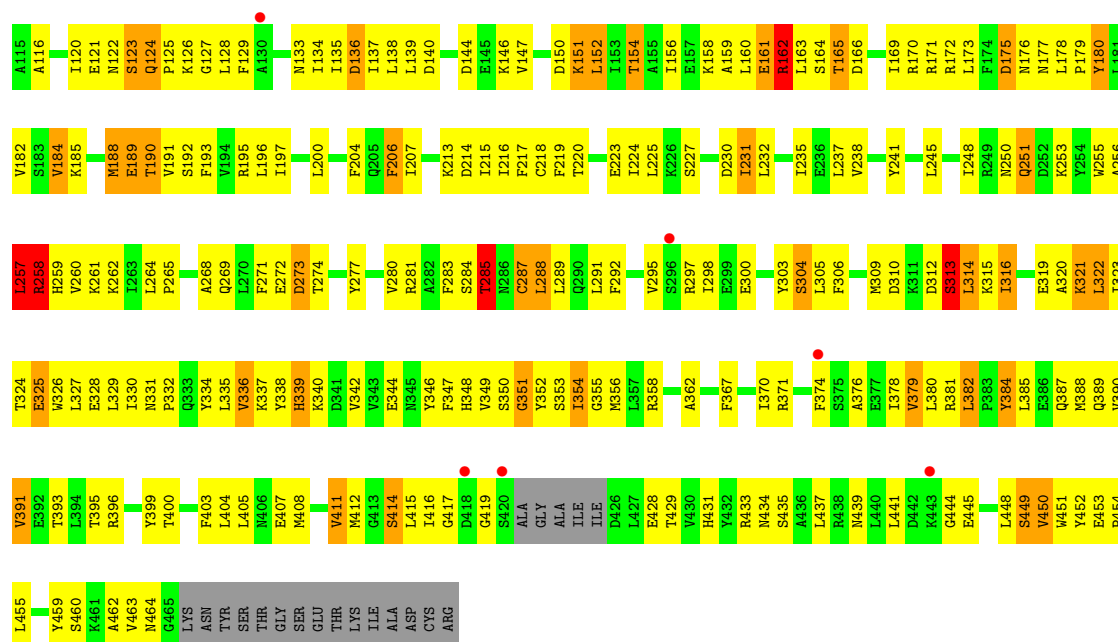
Chain	Residue	Modelled	Actual	Comment	Reference
C	-14	HIS	-	EXPRESSION TAG	UNP P38348
C	-13	HIS	-	EXPRESSION TAG	UNP P38348
C	-12	HIS	-	EXPRESSION TAG	UNP P38348
C	-11	HIS	-	EXPRESSION TAG	UNP P38348
C	-10	HIS	-	EXPRESSION TAG	UNP P38348
C	-9	SER	-	EXPRESSION TAG	UNP P38348
C	-8	SER	-	EXPRESSION TAG	UNP P38348
C	-7	GLY	-	EXPRESSION TAG	UNP P38348
C	-6	LEU	-	EXPRESSION TAG	UNP P38348
C	-5	VAL	-	EXPRESSION TAG	UNP P38348
C	-4	PRO	-	EXPRESSION TAG	UNP P38348
C	-3	ARG	-	EXPRESSION TAG	UNP P38348
C	-2	GLY	-	EXPRESSION TAG	UNP P38348
C	-1	SER	-	EXPRESSION TAG	UNP P38348
C	0	HIS	-	EXPRESSION TAG	UNP P38348

- Molecule 2 is a protein called 26S protease regulatory subunit 7 homolog.

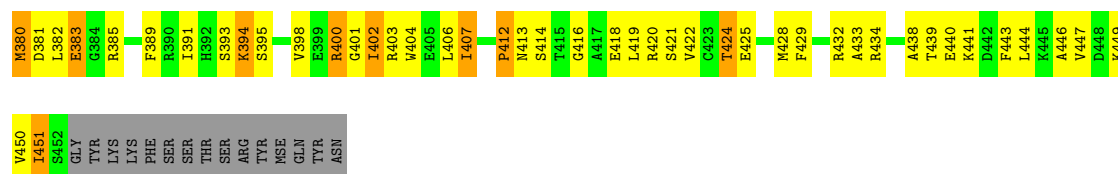
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	73	Total	C	N	O	S	Se	0	0	0
			575	357	110	103	2	3			
2	D	73	Total	C	N	O	S	Se	0	0	0
			575	357	110	103	2	3			

There are 2 discrepancies between the modelled and reference sequences:

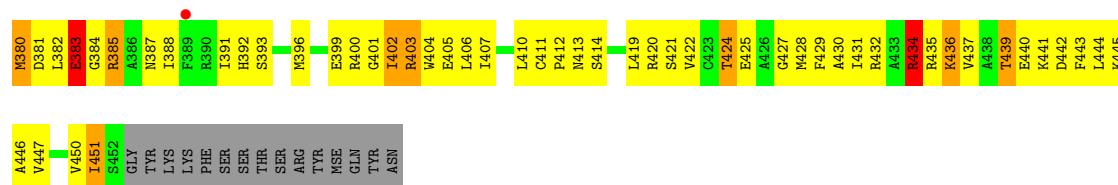
Chain	Residue	Modelled	Actual	Comment	Reference
B	380	MSE	-	EXPRESSION TAG	UNP P33299
D	380	MSE	-	EXPRESSION TAG	UNP P33299



- Molecule 2: 26S protease regulatory subunit 7 homolog



- Molecule 2: 26S protease regulatory subunit 7 homolog



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	187.28Å 187.28Å 379.57Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 3.80 46.82 – 3.79	Depositor EDS
% Data completeness (in resolution range)	(Not available) (30.00-3.80) 99.1 (46.82-3.79)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.80 (at 3.77Å)	Xtriage
Refinement program	CNS 1.3	Depositor
R, R_{free}	0.251 , 0.278 0.251 , 0.277	Depositor DCC
R_{free} test set	1952 reflections (4.97%)	DCC
Wilson B-factor (Å ²)	143.2	Xtriage
Anisotropy	0.147	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 129.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8594	wwPDB-VP
Average B, all atoms (Å ²)	163.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.48% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

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	0.38	0/3785	0.47	0/5114
1	C	0.36	0/3775	0.49	0/5099
2	B	0.34	0/579	0.47	0/769
2	D	0.36	0/579	0.47	0/769
All	All	0.37	0/8718	0.48	0/11751

There are no bond length outliers.

There are no bond angle outliers.

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	3727	0	3763	236	0
1	C	3717	0	3742	309	0
2	B	575	0	598	50	0
2	D	575	0	598	65	0
All	All	8594	0	8701	642	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 37.

The worst 5 of 642 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:A:464:ASN:OD1	1:C:82:LYS:NZ	1.93	1.01
1:C:88:ASP:HB3	1:C:91:ASP:HB2	1.42	1.01
1:C:124:GLN:HB2	1:C:125:PRO:HD3	1.43	0.98
1:A:320:ALA:HA	1:A:323:ILE:HG13	1.44	0.98
1:A:321:LYS:HD2	1:A:321:LYS:H	1.34	0.90

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	455/500 (91%)	366 (80%)	67 (15%)	22 (5%)	2	30
1	C	450/500 (90%)	330 (73%)	83 (18%)	37 (8%)	1	17
2	B	71/88 (81%)	55 (78%)	11 (16%)	5 (7%)	1	22
2	D	71/88 (81%)	50 (70%)	15 (21%)	6 (8%)	1	16
All	All	1047/1176 (89%)	801 (76%)	176 (17%)	70 (7%)	1	23

5 of 70 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	124	GLN
1	A	190	THR
1	A	231	ILE
1	A	281	ARG
1	A	449	SER

5.3.2 Protein sidechains [i](#)

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	424/452 (94%)	385 (91%)	39 (9%)	11	43
1	C	424/452 (94%)	382 (90%)	42 (10%)	9	40
2	B	62/72 (86%)	55 (89%)	7 (11%)	7	34
2	D	62/72 (86%)	53 (86%)	9 (14%)	4	24
All	All	972/1048 (93%)	875 (90%)	97 (10%)	9	39

5 of 97 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	424	THR
1	C	91	ASP
2	D	383	GLU
2	B	451	ILE
1	C	19	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 23 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	434	ASN
1	C	122	ASN
1	C	389	GLN
1	C	24	ASN
1	C	124	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

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

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

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2	OWAB(Å ²)	Q < 0.9
1	A	449/500 (89%)	0.32	23 (5%) 29 23	87, 158, 255, 289	0
1	C	446/500 (89%)	0.21	12 (2%) 55 44	102, 156, 216, 274	0
2	B	70/88 (79%)	0.09	0 100 100	106, 154, 195, 221	0
2	D	70/88 (79%)	0.12	1 (1%) 75 67	98, 144, 178, 202	0
All	All	1035/1176 (88%)	0.25	36 (3%) 44 35	87, 156, 236, 289	0

The worst 5 of 36 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	70	SER	6.5
1	A	13	LEU	4.6
1	A	29	ILE	4.5
1	A	421	ALA	4.5
1	A	422	GLY	4.1

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.