Full wwPDB X-ray Structure Validation Report (i)

Jan 2, 2024 – 12:17 pm GMT

PDB ID	:	5FWI
Title	:	structure of usp7 catalytic domain and three ubl-domains
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Deposited on	:	2016-02-17
Resolution	:	3.40 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

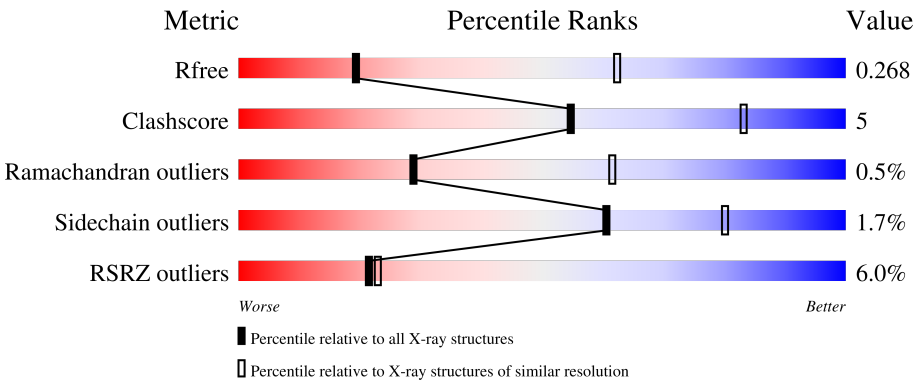
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.40 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}	130704	1026 (3.48-3.32)
Clashscore	141614	1055 (3.48-3.32)
Ramachandran outliers	138981	1038 (3.48-3.32)
Sidechain outliers	138945	1038 (3.48-3.32)
RSRZ outliers	127900	2173 (3.50-3.30)

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 of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. 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 <=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	C	680	6% 86% 11% • •

2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 5352 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
			Total	C	N	O	S			
1	C	662	5352	3394	904	1021	33	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

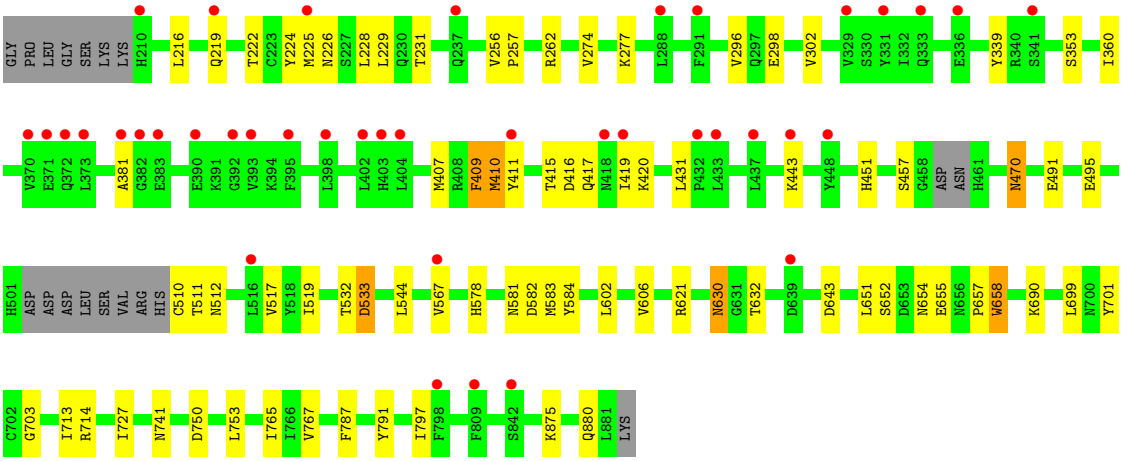
Chain	Residue	Modelled	Actual	Comment	Reference
C	203	GLY	-	expression tag	UNP Q93009
C	204	PRO	-	expression tag	UNP Q93009
C	205	LEU	-	expression tag	UNP Q93009
C	206	GLY	-	expression tag	UNP Q93009

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: UBIQUITIN CARBOXYL-TERMINAL HYDROLASE 7

Chain C:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	F 2 2 2	Depositor
Cell constants a, b, c, α, β, γ	115.18Å 195.04Å 219.78Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	109.89 – 3.40 48.76 – 3.40	Depositor EDS
% Data completeness (in resolution range)	99.8 (109.89-3.40) 99.9 (48.76-3.40)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.71 (at 3.40Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.222 , 0.267 0.225 , 0.268	Depositor DCC
R_{free} test set	851 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å^2)	120.2	Xtriage
Anisotropy	0.525	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.28 , 106.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	5352	wwPDB-VP
Average B, all atoms (Å^2)	171.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.25% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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	C	0.44	0/5467	0.66	1/7386 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	621	ARG	NE-CZ-NH1	5.32	122.96	120.30

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	C	5352	0	5217	55	0
All	All	5352	0	5217	55	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 5.

All (55) 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:C:409:PHE:HB3	1:C:420:LYS:HA	1.60	0.83
1:C:225:MET:O	1:C:229:LEU:HG	1.89	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:224:TYR:CD1	1:C:296:VAL:HG22	2.30	0.67
1:C:219:GLN:O	1:C:222:THR:HG23	1.94	0.67
1:C:224:TYR:CG	1:C:296:VAL:HG22	2.32	0.65
1:C:630:ASN:HD22	1:C:630:ASN:C	2.03	0.62
1:C:216:LEU:HB2	1:C:226:ASN:ND2	2.14	0.62
1:C:457:SER:HB3	1:C:511:THR:HG22	1.83	0.61
1:C:583:MET:CE	1:C:765:ILE:HD11	2.31	0.61
1:C:419:ILE:HG13	1:C:420:LYS:H	1.67	0.58
1:C:643:ASP:HB3	1:C:651:LEU:HD21	1.84	0.57
1:C:360:ILE:HD11	1:C:431:LEU:HD13	1.87	0.57
1:C:652:SER:O	1:C:655:GLU:HG3	2.05	0.56
1:C:652:SER:HA	1:C:658:TRP:CZ2	2.40	0.56
1:C:353:SER:HA	1:C:407:MET:HG2	1.87	0.55
1:C:411:TYR:HA	1:C:417:GLN:O	2.07	0.55
1:C:602:LEU:O	1:C:606:VAL:HG23	2.07	0.55
1:C:510:CYS:SG	1:C:511:THR:N	2.79	0.52
1:C:228:LEU:HD12	1:C:228:LEU:O	2.10	0.52
1:C:690:LYS:HB2	1:C:767:VAL:HG22	1.92	0.51
1:C:407:MET:SD	1:C:407:MET:N	2.85	0.50
1:C:339:TYR:CD2	1:C:381:ALA:HB1	2.47	0.49
1:C:583:MET:HE2	1:C:765:ILE:HD11	1.94	0.49
1:C:630:ASN:ND2	1:C:632:THR:OG1	2.41	0.48
1:C:658:TRP:HE3	1:C:658:TRP:HA	1.79	0.48
1:C:224:TYR:CE1	1:C:296:VAL:HA	2.48	0.48
1:C:532:THR:HG22	1:C:533:ASP:N	2.28	0.48
1:C:658:TRP:HA	1:C:658:TRP:CE3	2.48	0.47
1:C:298:GLU:O	1:C:302:VAL:HG23	2.14	0.47
1:C:567:VAL:HG22	1:C:658:TRP:HB3	1.96	0.47
1:C:470:ASN:HD22	1:C:470:ASN:H	1.63	0.46
1:C:419:ILE:CG1	1:C:420:LYS:H	2.28	0.46
1:C:231:THR:HG21	1:C:517:VAL:HG21	1.97	0.46
1:C:262:ARG:HG3	1:C:544:LEU:HD22	1.98	0.45
1:C:727:ILE:HD12	1:C:727:ILE:H	1.82	0.45
1:C:714:ARG:NH1	1:C:750:ASP:OD1	2.50	0.44
1:C:216:LEU:CB	1:C:226:ASN:ND2	2.81	0.44
1:C:410:MET:O	1:C:419:ILE:HG22	2.17	0.44
1:C:581:ASN:O	1:C:582:ASP:HB2	2.18	0.43
1:C:699:LEU:HD11	1:C:787:PHE:CZ	2.54	0.43
1:C:713:ILE:HD13	1:C:753:LEU:HD23	1.99	0.43
1:C:419:ILE:HG13	1:C:420:LYS:N	2.31	0.43
1:C:219:GLN:HE22	1:C:277:LYS:HG3	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:491:GLU:HG3	1:C:495:GLU:HG3	2.01	0.42
1:C:451:HIS:CD2	1:C:519:ILE:HD12	2.53	0.42
1:C:584:TYR:C	1:C:584:TYR:CD1	2.93	0.41
1:C:741:ASN:HB2	1:C:791:TYR:CG	2.56	0.41
1:C:690:LYS:O	1:C:767:VAL:HA	2.20	0.41
1:C:701:TYR:CE2	1:C:703:GLY:HA2	2.55	0.41
1:C:630:ASN:C	1:C:630:ASN:ND2	2.73	0.41
1:C:578:HIS:ND1	1:C:584:TYR:HA	2.35	0.41
1:C:216:LEU:HD23	1:C:274:VAL:HB	2.02	0.40
1:C:256:VAL:N	1:C:257:PRO:CD	2.85	0.40
1:C:578:HIS:CE1	1:C:584:TYR:C	2.95	0.40
1:C:797:ILE:O	1:C:875:LYS:HA	2.22	0.40

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	C	656/680 (96%)	634 (97%)	19 (3%)	3 (0%)	29 61

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	415	THR
1	C	416	ASP
1	C	657	PRO

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	C	591/617 (96%)	581 (98%)	10 (2%)	60 80

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	409	PHE
1	C	410	MET
1	C	443	LYS
1	C	470	ASN
1	C	512	ASN
1	C	533	ASP
1	C	630	ASN
1	C	654	ASN
1	C	658	TRP
1	C	880	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	387	GLN
1	C	470	ASN
1	C	617	GLN
1	C	630	ASN
1	C	654	ASN
1	C	838	GLN
1	C	880	GLN

5.3.3 RNA (i)

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 monosaccharides 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)

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	C	662/680 (97%)	0.28	40 (6%)	21	23	86, 166, 268, 352	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	372	GLN	7.3
1	C	370	VAL	5.0
1	C	433	LEU	4.6
1	C	403	HIS	4.4
1	C	419	ILE	3.9
1	C	373	LEU	3.8
1	C	331	TYR	3.6
1	C	398	LEU	3.6
1	C	219	GLN	3.5
1	C	291	PHE	3.5
1	C	437	LEU	3.4
1	C	288	LEU	3.3
1	C	393	VAL	3.1
1	C	237	GLN	3.1
1	C	392	GLY	3.1
1	C	382	GLY	2.9
1	C	395	PHE	2.9
1	C	443	LYS	2.8
1	C	411	TYR	2.8
1	C	432	PRO	2.8
1	C	404	LEU	2.8
1	C	390	GLU	2.7
1	C	371	GLU	2.7
1	C	516	LEU	2.7
1	C	639	ASP	2.7
1	C	798	PHE	2.7
1	C	381	ALA	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	333	GLN	2.6
1	C	448	TYR	2.6
1	C	418	ASN	2.5
1	C	567	VAL	2.4
1	C	341	SER	2.3
1	C	225	MET	2.3
1	C	336	GLU	2.2
1	C	402	LEU	2.1
1	C	383	GLU	2.1
1	C	842	SER	2.1
1	C	809	PHE	2.0
1	C	210	HIS	2.0
1	C	329	VAL	2.0

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 monosaccharides 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.