



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:49 AM GMT

PDB ID : 3Q5X
Title : Structure of proteasome tether
Authors : Schumacher, M.A.
Deposited on : 2010-12-30
Resolution : 2.98 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
<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
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

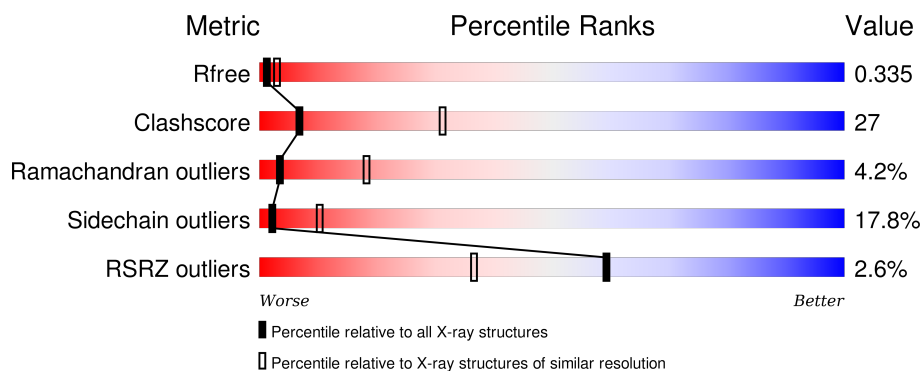
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.98 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}	91344	1992 (3.00-2.96)
Clashscore	102246	2349 (3.00-2.96)
Ramachandran outliers	100387	2274 (3.00-2.96)
Sidechain outliers	100360	2277 (3.00-2.96)
RSRZ outliers	91569	2007 (3.00-2.96)

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	245	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1570 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 Protein cut8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	193	Total	C	N	O	S	0	0	0
			1569	1003	270	288	8			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	EXPRESSION TAG	UNP P38937
A	-18	GLY	-	EXPRESSION TAG	UNP P38937
A	-17	SER	-	EXPRESSION TAG	UNP P38937
A	-16	SER	-	EXPRESSION TAG	UNP P38937
A	-15	HIS	-	EXPRESSION TAG	UNP P38937
A	-14	HIS	-	EXPRESSION TAG	UNP P38937
A	-13	HIS	-	EXPRESSION TAG	UNP P38937
A	-12	HIS	-	EXPRESSION TAG	UNP P38937
A	-11	HIS	-	EXPRESSION TAG	UNP P38937
A	-10	HIS	-	EXPRESSION TAG	UNP P38937
A	-9	SER	-	EXPRESSION TAG	UNP P38937
A	-8	SER	-	EXPRESSION TAG	UNP P38937
A	-7	GLY	-	EXPRESSION TAG	UNP P38937
A	-6	LEU	-	EXPRESSION TAG	UNP P38937
A	-5	VAL	-	EXPRESSION TAG	UNP P38937
A	-4	PRO	-	EXPRESSION TAG	UNP P38937
A	-3	ARG	-	EXPRESSION TAG	UNP P38937
A	-2	GLY	-	EXPRESSION TAG	UNP P38937
A	-1	SER	-	EXPRESSION TAG	UNP P38937
A	0	HIS	-	EXPRESSION TAG	UNP P38937

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	O	0	0
			1	1		

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	143.44Å 36.20Å 53.99Å 90.00° 108.17° 90.00°	Depositor
Resolution (Å)	34.99 – 2.98 34.99 – 2.69	Depositor EDS
% Data completeness (in resolution range)	95.7 (34.99-2.98) 95.6 (34.99-2.69)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.93 (at 2.68Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.262 , 0.295 0.292 , 0.335	Depositor DCC
R_{free} test set	738 reflections (13.86%)	DCC
Wilson B-factor (Å ²)	66.9	Xtriage
Anisotropy	0.730	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 42.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.53$, $\langle L^2 \rangle = 0.37$	Xtriage
Outliers	0 of 7244 reflections	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	1570	wwPDB-VP
Average B, all atoms (Å ²)	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.92% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.79	5/1606 (0.3%)	0.77	3/2184 (0.1%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	196	GLU	CD-OE1	-6.90	1.18	1.25
1	A	196	GLU	CD-OE2	-6.89	1.18	1.25
1	A	56	GLU	CD-OE1	-5.46	1.19	1.25
1	A	112	MET	CG-SD	5.35	1.95	1.81
1	A	56	GLU	CG-CD	-5.14	1.44	1.51

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	127	PRO	CA-N-CD	-7.88	100.47	111.50
1	A	141	ASP	CB-CG-OD1	7.30	124.87	118.30
1	A	85	LEU	CB-CA-C	6.13	121.84	110.20

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	1569	0	1561	84	0
2	A	1	0	0	1	0
All	All	1570	0	1561	84	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 27.

All (84) 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:127:PRO:CG	1:A:175:VAL:HG12	1.55	1.36
1:A:127:PRO:HG2	1:A:175:VAL:CG1	1.52	1.36
1:A:132:CYS:SG	1:A:135:LYS:CD	2.17	1.33
1:A:132:CYS:SG	1:A:135:LYS:HD3	1.68	1.32
1:A:132:CYS:SG	1:A:135:LYS:HE2	1.88	1.12
1:A:132:CYS:SG	1:A:135:LYS:CE	2.39	1.10
1:A:85:LEU:HD11	1:A:149:GLU:HB3	1.42	0.98
1:A:35:LEU:N	2:A:226:HOH:O	1.95	0.96
1:A:132:CYS:HG	1:A:135:LYS:HD3	1.12	0.93
1:A:175:VAL:HG13	1:A:178:ARG:HH21	1.35	0.91
1:A:106:ARG:HB3	1:A:106:ARG:HH11	1.36	0.89
1:A:132:CYS:SG	1:A:135:LYS:CG	2.62	0.88
1:A:127:PRO:HG2	1:A:175:VAL:HG12	0.87	0.86
1:A:127:PRO:HG3	1:A:175:VAL:HG12	1.62	0.81
1:A:127:PRO:HG2	1:A:175:VAL:HG11	1.61	0.75
1:A:164:GLN:O	1:A:168:GLU:HG2	1.88	0.74
1:A:33:LEU:H	1:A:33:LEU:HD23	1.51	0.74
1:A:142:ALA:O	1:A:146:VAL:HG23	1.89	0.71
1:A:175:VAL:HG13	1:A:178:ARG:NH2	2.04	0.71
1:A:88:LEU:HD12	1:A:88:LEU:C	2.15	0.67
1:A:108:ARG:O	1:A:112:MET:HG2	1.99	0.63
1:A:198:ASN:HD21	1:A:203:ASN:HD22	1.46	0.63
1:A:134:GLU:CD	1:A:134:GLU:H	2.03	0.62
1:A:185:VAL:O	1:A:187:PRO:HD3	2.00	0.61
1:A:123:PRO:HA	1:A:126:LEU:HD12	1.83	0.61
1:A:133:PHE:CE2	1:A:180:LEU:HD22	2.36	0.61
1:A:201:SER:O	1:A:204:ARG:HD3	1.99	0.60
1:A:88:LEU:HD13	1:A:107:ILE:HG21	1.84	0.59
1:A:198:ASN:HD21	1:A:203:ASN:ND2	2.01	0.58
1:A:150:LEU:O	1:A:162:LYS:NZ	2.36	0.58
1:A:189:LEU:HD21	1:A:215:LEU:HD13	1.85	0.58
1:A:106:ARG:CB	1:A:106:ARG:HH11	2.13	0.57
1:A:133:PHE:CZ	1:A:180:LEU:HD22	2.40	0.56
1:A:130:SER:OG	1:A:136:ASN:OD1	2.22	0.55
1:A:173:TRP:O	1:A:177:LEU:HG	2.07	0.54
1:A:132:CYS:SG	1:A:135:LYS:HG2	2.46	0.54
1:A:171:GLY:O	1:A:175:VAL:HG23	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:132:CYS:SG	1:A:135:LYS:HB2	2.49	0.53
1:A:123:PRO:HA	1:A:126:LEU:CD1	2.39	0.52
1:A:63:ARG:CZ	1:A:63:ARG:HB2	2.38	0.52
1:A:110:LYS:O	1:A:114:VAL:HG23	2.09	0.52
1:A:88:LEU:O	1:A:88:LEU:HD12	2.10	0.52
1:A:121:MET:O	1:A:124:CYS:HB2	2.10	0.52
1:A:92:PHE:HB3	1:A:93:PRO:HD2	1.91	0.51
1:A:82:ARG:O	1:A:86:ILE:HG13	2.10	0.51
1:A:180:LEU:HD11	1:A:187:PRO:HD2	1.92	0.51
1:A:198:ASN:ND2	1:A:203:ASN:HD22	2.09	0.51
1:A:47:LEU:O	1:A:51:LEU:HB2	2.11	0.50
1:A:48:PHE:O	1:A:52:LEU:HD23	2.12	0.50
1:A:88:LEU:HD23	1:A:114:VAL:HG21	1.93	0.50
1:A:122:VAL:N	1:A:123:PRO:HD2	2.27	0.50
1:A:127:PRO:HB2	1:A:128:PRO:HD3	1.94	0.49
1:A:191:LEU:HA	1:A:212:LEU:HD11	1.94	0.49
1:A:46:GLN:O	1:A:50:ILE:HG23	2.13	0.49
1:A:166:TYR:CE2	1:A:201:SER:HB2	2.46	0.49
1:A:107:ILE:O	1:A:108:ARG:C	2.50	0.49
1:A:133:PHE:CE2	1:A:137:ILE:HD11	2.49	0.48
1:A:132:CYS:SG	1:A:135:LYS:CB	3.02	0.48
1:A:88:LEU:HD13	1:A:107:ILE:CG2	2.44	0.47
1:A:33:LEU:HD12	1:A:37:ARG:HG3	1.98	0.46
1:A:176:VAL:O	1:A:179:GLN:HG2	2.15	0.46
1:A:118:LEU:O	1:A:122:VAL:HG23	2.15	0.46
1:A:92:PHE:HA	1:A:107:ILE:HD11	1.96	0.46
1:A:108:ARG:HE	1:A:112:MET:HG3	1.81	0.46
1:A:145:ASN:HA	1:A:148:HIS:CD2	2.51	0.46
1:A:187:PRO:O	1:A:189:LEU:HD13	2.15	0.46
1:A:182:ASP:O	1:A:184:PRO:HD3	2.16	0.46
1:A:92:PHE:CD2	1:A:98:LYS:HB3	2.52	0.45
1:A:77:CYS:O	1:A:80:THR:HB	2.15	0.45
1:A:180:LEU:CD2	1:A:189:LEU:HD11	2.48	0.43
1:A:155:ASN:OD1	1:A:156:PRO:HD2	2.19	0.43
1:A:189:LEU:HB3	1:A:212:LEU:HD21	2.01	0.43
1:A:57:LYS:HB3	1:A:58:HIS:CD2	2.54	0.43
1:A:97:ASP:C	1:A:99:ARG:H	2.23	0.43
1:A:51:LEU:O	1:A:54:CYS:HB2	2.18	0.42
1:A:133:PHE:CD2	1:A:137:ILE:HD11	2.54	0.42
1:A:186:VAL:HG11	1:A:215:LEU:O	2.20	0.41
1:A:213:LYS:O	1:A:216:GLN:HB2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:189:LEU:HA	1:A:190:PRO:HD3	1.89	0.41
1:A:88:LEU:C	1:A:88:LEU:CD1	2.85	0.41
1:A:212:LEU:HD13	1:A:212:LEU:C	2.40	0.41
1:A:189:LEU:HD23	1:A:212:LEU:HD22	2.02	0.41
1:A:41:TYR:CD1	1:A:41:TYR:N	2.89	0.40
1:A:71:ALA:HB1	1:A:72:PRO:HD2	2.03	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	A	191/245 (78%)	166 (87%)	17 (9%)	8 (4%)	3	18

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	181	GLU
1	A	94	TYR
1	A	127	PRO
1	A	95	GLY
1	A	98	LYS
1	A	185	VAL
1	A	59	PRO
1	A	128	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	A	174/224 (78%)	143 (82%)	31 (18%)	2	10

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	ARG
1	A	35	LEU
1	A	36	SER
1	A	42	SER
1	A	44	LYS
1	A	51	LEU
1	A	56	GLU
1	A	66	ARG
1	A	74	MET
1	A	79	GLU
1	A	85	LEU
1	A	88	LEU
1	A	94	TYR
1	A	106	ARG
1	A	108	ARG
1	A	109	GLU
1	A	110	LYS
1	A	111	TYR
1	A	115	LEU
1	A	118	LEU
1	A	138	THR
1	A	149	GLU
1	A	163	SER
1	A	183	ARG
1	A	189	LEU
1	A	200	THR
1	A	204	ARG
1	A	207	GLU
1	A	209	LEU
1	A	214	GLN
1	A	216	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	26	GLN
1	A	32	GLN
1	A	145	ASN
1	A	203	ASN
1	A	210	ASN

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 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](#)

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	193/245 (78%)	0.37	5 (2%) 59 37	46, 70, 98, 110	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	186	VAL	4.2
1	A	180	LEU	3.4
1	A	94	TYR	3.3
1	A	184	PRO	2.8
1	A	188	LEU	2.6

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.