

wwPDB EM Validation Summary Report (i)

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PDB ID	:	5VGZ
EMDB ID	:	EMD-8672
Title	:	Conformational Landscape of the p28-Bound Human Proteasome Regulatory
		Particle
Authors	:	Lu, Y.; Wu, J.; Dong, Y.; Chen, S.; Sun, S.; Ma, Y.B.; Ouyang, Q.; Finley,
		D.; Kirschner, M.W.; Mao, Y.
Deposited on	:	2017-04-12
Resolution	:	4.50 Å(reported)

This is a wwPDB EM Validation Summary 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/EMValidationReportHelp 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:

EMDB validation analysis	:	0.0.1. dev 50
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ	:	1.9.9
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $ELECTRON\ MICROSCOPY$

The reported resolution of this entry is 4.50 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)	${f EM} {f structures} \ (\#{f Entries})$
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for $\geq=3, 2, 1$ 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 $\leq=5\%$ The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion < 40%). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain	
1	А	83	99%	•
2	В	73	14%	
3	С	118	9%	·
4	D	107	9%	·
5	Е	104	26%	
6	F	115	85%	• 12%
7	U	935	90%	• 10%
8	V	488	28%	•
9	W	456	98%	·

Continued on next page...



Mal	Chain	Longth	Quality of choir
IVIOI	Unain	Length	Quality of chain
			25%
10	Х	385	100%
			5%
11	Y	378	100%
			7%
12	Z	286	97% •
			6%
13	a	374	100%
			7%
14	b	191	99% •
			8%
15	с	287	99%
			21%
16	d	257	100%
			39%
17	е	70	97% •

Continued from previous page...



2 Entry composition (i)

There are 18 unique types of molecules in this entry. The entry contains 36741 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, 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 26S proteasome regulatory subunit 7.

Mol	Chain	Residues		At	oms	AltConf	Trace		
1	А	83	Total 654	C 416	N 111	0 125	${S \over 2}$	0	0

• Molecule 2 is a protein called 26S proteasome regulatory subunit 4.

Mol	Chain	Residues		Ate	\mathbf{oms}	AltConf	Trace		
2	В	73	Total 556	C 346	N 92	0 115	${ m S} { m 3}$	0	0

• Molecule 3 is a protein called 26S proteasome regulatory subunit 8.

Mol	Chain	Residues		At	oms	AltConf	Trace		
3	С	118	Total 954	C 597	N 176	0 179	${ m S} { m 2}$	0	0

• Molecule 4 is a protein called 26S proteasome regulatory subunit 6B.

Mol	Chain	Residues		Ato	\mathbf{ms}	AltConf	Trace	
4	D	107	Total 870	C 555	N 147	O 168	0	0

• Molecule 5 is a protein called 26S proteasome regulatory subunit 10B.

Mol	Chain	Residues		At	oms	AltConf	Trace		
5	Е	104	Total 853	C 537	N 155	0 158	${ m S} { m 3}$	0	0

• Molecule 6 is a protein called 26S proteasome regulatory subunit 6A.

Mol	Chain	Residues		At	oms	AltConf	Trace		
6	F	101	Total 799	C 512	N 133	0 151	${ m S} { m 3}$	0	0



• Molecule 7 is a protein called 26S proteasome non-ATPase regulatory subunit 1.

Mol	Chain	Residues		Α	AltConf	Trace			
7	U	845	Total 6568	C 4169	N 1118	O 1236	$\begin{array}{c} \mathrm{S} \\ 45 \end{array}$	0	0

• Molecule 8 is a protein called 26S proteasome non-ATPase regulatory subunit 3.

Mol	Chain	Residues	Atoms			AltConf	Trace		
8	V	488	Total 3919	C 2487	N 695	0 723	S 14	0	0

• Molecule 9 is a protein called 26S proteasome non-ATPase regulatory subunit 12.

Mol	Chain	Residues	Atoms				AltConf	Trace	
9	W	456	Total 3703	C 2339	N 635	0 704	S 25	0	0

• Molecule 10 is a protein called 26S proteasome non-ATPase regulatory subunit 11.

Mol	Chain	Residues	Atoms				AltConf	Trace	
10	Х	385	Total 3048	C 1939	N 515	0 582	S 12	0	0

• Molecule 11 is a protein called 26S proteasome non-ATPase regulatory subunit 6.

Mol	Chain	Residues	Atoms				AltConf	Trace	
11	Y	378	Total 3115	C 1987	N 533	0 578	S 17	0	0

• Molecule 12 is a protein called 26S proteasome non-ATPase regulatory subunit 7.

Mol	Chain	Residues	Atoms				AltConf	Trace	
12	Z	286	Total 2281	C 1457	N 392	0 427	${ m S}{ m 5}$	0	0

• Molecule 13 is a protein called 26S proteasome non-ATPase regulatory subunit 13.

Mol	Chain	Residues	Atoms			AltConf	Trace		
13	a	374	Total 3003	C 1915	N 511	O 562	S 15	0	0

• Molecule 14 is a protein called 26S proteasome non-ATPase regulatory subunit 4.



Mol	Chain	Residues	Atoms			AltConf	Trace		
14	b	191	Total 1458	C 910	N 261	O 279	S 8	0	0

• Molecule 15 is a protein called 26S proteasome non-ATPase regulatory subunit 14.

Mol	Chain	Residues	Atoms				AltConf	Trace	
15	с	287	Total 2260	C 1430	N 389	O 422	S 19	0	0

• Molecule 16 is a protein called 26S proteasome non-ATPase regulatory subunit 8.

Mol	Chain	Residues	Atoms				AltConf	Trace	
16	d	257	Total 2116	C 1371	N 346	O 390	S 9	0	0

• Molecule 17 is a protein called 26S proteasome complex subunit SEM1.

Mol	Chain	Residues	Atoms				AltConf	Trace	
17	е	70	Total 583	C 357	N 89	0 135	${ m S} { m 2}$	0	0

• Molecule 18 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	AltConf
18	с	1	Total Zn 1 1	0



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 atom inclusion in map 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 diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). 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: 26S proteasome regulatory subunit 7





• Molecule 6: 26S proteasome regulatory subunit 6A 18% Chain F: 85% 12% • Molecule 7: 26S proteasome non-ATPase regulatory subunit 1 Chain U: 90% 10% ILE ALA SER VAL SER SER ALA PHE VAL VAL CYS LYS LYS LYS THR PRO GLU GLU SER SER PRO LYS GLU CYS GLU CYS GLU LYS VAL VAL SER THR • Molecule 8: 26S proteasome non-ATPase regulatory subunit 3 28% Chain V: 96% E83 K84 • Molecule 9: 26S proteasome non-ATPase regulatory subunit 12











4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	117471	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE	Depositor
	CORRECTION	
Microscope	FEI TECNAI ARCTICA	Depositor
Voltage (kV)	200	Depositor
Electron dose $(e^-/\text{\AA}^2)$	50	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.015	Depositor
Minimum map value	-0.005	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.001	Depositor
Recommended contour level	0.004	Depositor
Map size (Å)	250.88, 250.88, 250.88	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles $(^{\circ})$	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.98, 0.98, 0.98	Depositor



5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN

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).

Mal	Chain	Bond	lengths	B	ond angles
	Ullaili	RMSZ	# Z > 5	RMSZ	# Z > 5
1	А	0.32	0/668	0.66	1/911~(0.1%)
2	В	0.33	0/563	0.66	0/763
3	С	0.30	0/963	0.61	0/1294
4	D	0.32	0/884	0.63	1/1195~(0.1%)
5	Е	0.29	0/861	0.56	0/1149
6	F	0.29	0/807	0.66	3/1086~(0.3%)
7	U	0.31	0/6684	0.59	2/9043~(0.0%)
8	V	0.32	0/3998	0.68	3/5402~(0.1%)
9	W	0.31	0/3750	0.60	3/5039~(0.1%)
10	Х	0.29	0/3091	0.51	0/4165
11	Y	0.31	0/3173	0.57	1/4273~(0.0%)
12	Ζ	0.31	0/2324	0.60	1/3150~(0.0%)
13	a	0.31	0/3061	0.59	0/4144
14	b	0.30	0/1478	0.57	0/2001
15	с	0.32	0/2302	0.57	0/3110
16	d	0.30	0/2162	0.57	1/2919~(0.0%)
17	е	0.31	0/596	0.62	1/805~(0.1%)
All	All	0.31	0/37365	0.60	17/50449~(0.0%)

There are no bond length outliers.

The worst 5 of 17 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	$Observed(^{o})$	$Ideal(^{o})$
17	е	56	LEU	CA-CB-CG	6.89	131.15	115.30
16	d	122	LEU	CA-CB-CG	6.72	130.76	115.30
7	U	252	LEU	CA-CB-CG	6.59	130.46	115.30
8	V	170	LEU	CA-CB-CG	6.39	130.00	115.30
9	W	219	THR	O-C-N	5.92	132.18	122.70

There are no chirality outliers.



There are no planarity outliers.

5.2 Too-close contacts (i)

Due to software issues we are unable to calculate clashes - this section is therefore empty.

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 PDB entries followed by that with respect to all EM entries.

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	Perce	ntiles
1	А	81/83~(98%)	73~(90%)	8 (10%)	0	100	100
2	В	71/73~(97%)	57~(80%)	14 (20%)	0	100	100
3	С	116/118 (98%)	110 (95%)	5 (4%)	1 (1%)	17	56
4	D	105/107~(98%)	102 (97%)	3 (3%)	0	100	100
5	Е	102/104~(98%)	97~(95%)	5 (5%)	0	100	100
6	F	97/115 (84%)	91 (94%)	6 (6%)	0	100	100
7	U	837/935~(90%)	790 (94%)	46 (6%)	1 (0%)	51	85
8	V	486/488 (100%)	410 (84%)	67 (14%)	9 (2%)	8	41
9	W	452/456~(99%)	410 (91%)	38 (8%)	4 (1%)	17	56
10	Х	381/385~(99%)	371 (97%)	10 (3%)	0	100	100
11	Y	376/378~(100%)	355~(94%)	21 (6%)	0	100	100
12	Z	284/286~(99%)	260 (92%)	21 (7%)	3 (1%)	14	52
13	a	372/374~(100%)	346 (93%)	25~(7%)	1 (0%)	41	76
14	b	189/191 (99%)	174 (92%)	15 (8%)	0	100	100
15	с	285/287~(99%)	259 (91%)	25 (9%)	1 (0%)	34	72
16	d	255/257~(99%)	230 (90%)	25 (10%)	0	100	100
17	e	68/70~(97%)	61 (90%)	7 (10%)	0	100	100
All	All	4557/4707 (97%)	4196 (92%)	341 (8%)	20 (0%)	38	72

5 of 20 Ramachandran outliers are listed below:



Mol	Chain	Res	Type
3	С	91	PRO
8	V	27	PRO
8	V	190	ASP
8	V	193	GLN
8	V	225	ASP

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent side chain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

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	Perce	ntiles
1	А	73/73~(100%)	73~(100%)	0	100	100
2	В	66/66~(100%)	66 (100%)	0	100	100
3	С	107/107~(100%)	107 (100%)	0	100	100
4	D	98/98~(100%)	98 (100%)	0	100	100
5	Е	95/95~(100%)	95 (100%)	0	100	100
6	F	94/106~(89%)	94 (100%)	0	100	100
7	U	717/798~(90%)	714 (100%)	3 (0%)	91	94
8	V	422/422 (100%)	414 (98%)	8 (2%)	57	75
9	W	416/416 (100%)	410 (99%)	6 (1%)	67	81
10	Х	331/331~(100%)	331 (100%)	0	100	100
11	Y	334/334~(100%)	334 (100%)	0	100	100
12	Z	257/257~(100%)	252 (98%)	5 (2%)	57	75
13	a	334/334~(100%)	334 (100%)	0	100	100
14	b	167/167~(100%)	166 (99%)	1 (1%)	86	92
15	с	252/252~(100%)	251 (100%)	1 (0%)	91	94
16	d	231/231 (100%)	231 (100%)	0	100	100
17	е	63/63~(100%)	62 (98%)	1 (2%)	62	79
All	All	4057/4150 (98%)	4032 (99%)	25 (1%)	86	92

 $5~{\rm of}~25$ residues with a non-rotameric side chain are listed below:



Mol	Chain	Res	Type
9	W	220	GLU
12	Ζ	28	LYS
17	е	22	PHE
9	W	417	ARG
12	Ζ	90	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 46 such sidechains are listed below:

Mol	Chain	Res	Type
11	Y	48	ASN
13	а	372	HIS
12	Ζ	196	HIS
13	a	35	HIS
14	b	29	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)

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
9	W	1
10	Х	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	W	205:ILE	С	206:SER	N	3.21
1	Х	311:ALA	С	312:GLU	N	3.21



6 Map visualisation (i)

This section contains visualisations of the EMDB entry EMD-8672. These allow visual inspection of the internal detail of the map and identification of artifacts.

No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections (i)

6.1.1 Primary map



The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

6.2.1 Primary map



X Index: 128





Z Index: 128

The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

6.3.1 Primary map



X Index: 109

Y Index: 179

Z Index: 139

The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) (i)

6.4.1 Primary map



The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.



6.5 Orthogonal surface views (i)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.004. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.6 Mask visualisation (i)

This section was not generated. No masks/segmentation were deposited.



7 Map analysis (i)

This section contains the results of statistical analysis of the map.

7.1 Map-value distribution (i)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.



7.2 Volume estimate (i)



The volume at the recommended contour level is 523 $\rm nm^3;$ this corresponds to an approximate mass of 472 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.



7.3 Rotationally averaged power spectrum (i)



*Reported resolution corresponds to spatial frequency of 0.222 ${\rm \AA^{-1}}$



8 Fourier-Shell correlation (i)

This section was not generated. No FSC curve or half-maps provided.



9 Map-model fit (i)

This section contains information regarding the fit between EMDB map EMD-8672 and PDB model 5VGZ. Per-residue inclusion information can be found in section 3 on page 7.

9.1 Map-model overlay (i)



The images above show the 3D surface view of the map at the recommended contour level 0.004 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.



9.2 Q-score mapped to coordinate model (i)



The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model (i)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.004).



9.4 Atom inclusion (i)



At the recommended contour level, 89% of all backbone atoms, 71% of all non-hydrogen atoms, are inside the map.



1.0

0.0 <0.0

9.5 Map-model fit summary (i)

The table lists the average atom inclusion at the recommended contour level (0.004) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7070	0.1950
А	0.6990	0.1900
В	0.6620	0.1810
С	0.7190	0.2040
D	0.6900	0.2120
Е	0.5930	0.2090
F	0.6080	0.2080
U	0.8110	0.2220
V	0.5810	0.1630
W	0.6780	0.1800
Х	0.6060	0.1790
Y	0.7920	0.1890
Z	0.7400	0.2040
a	0.7750	0.1950
b	0.7640	0.1930
С	0.7240	0.2160
d	0.6430	0.1900
е	0.5290	0.1680

