



wwPDB X-ray Structure Validation Summary Report ⓘ

Nov 11, 2023 – 07:00 pm GMT

PDB ID : 4YA3
Title : Yeast 20S proteasome beta2-H116N mutant in complex with Ac-PAE-ep
Authors : Huber, E.M.; Groll, M.
Deposited on : 2015-02-17
Resolution : 2.70 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.4, CSD as541be (2020)
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

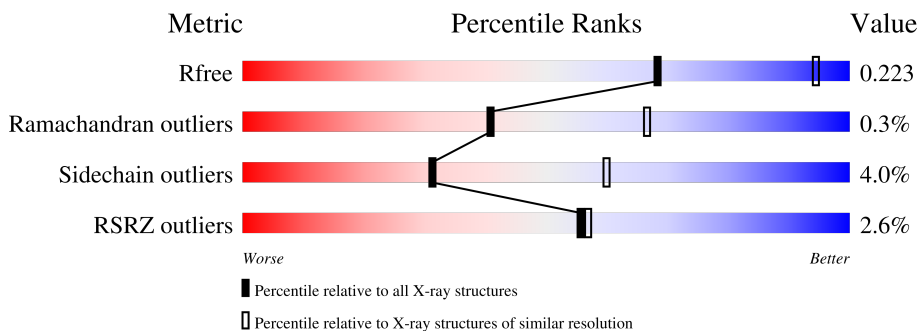
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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.



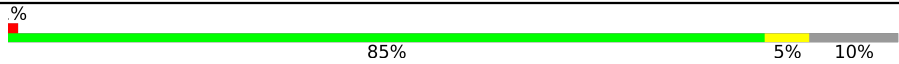

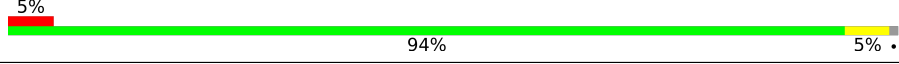
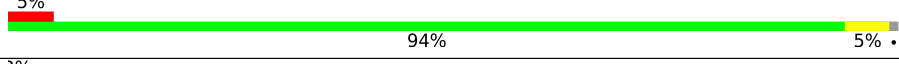

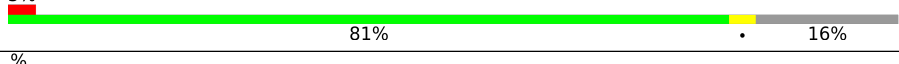
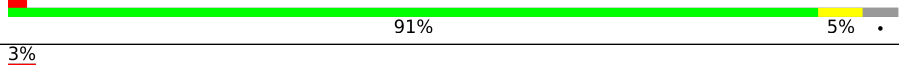
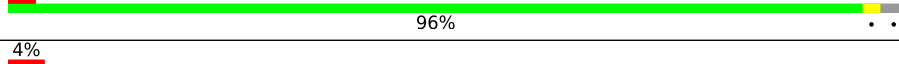
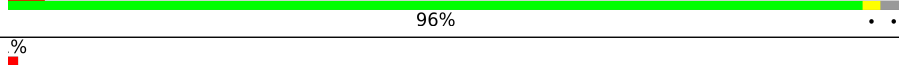
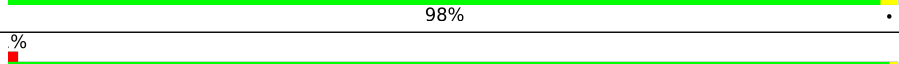
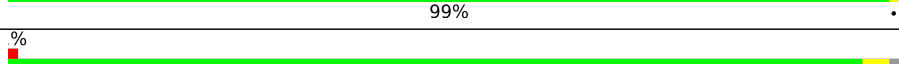
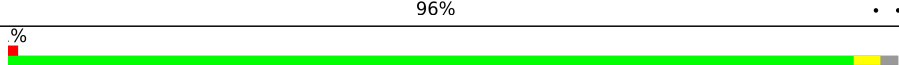
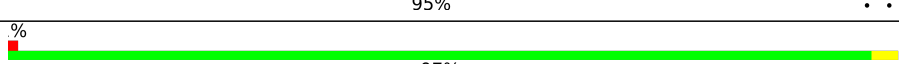
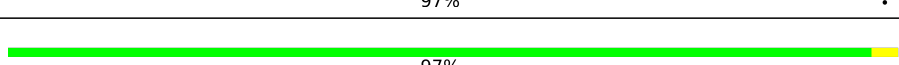
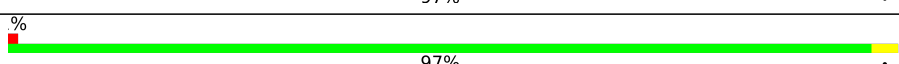
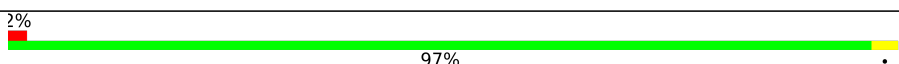
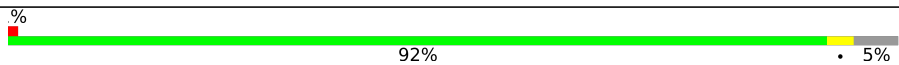
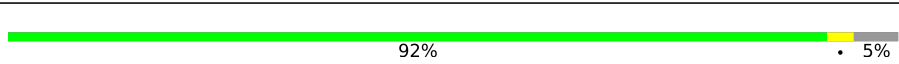
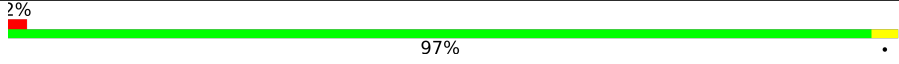
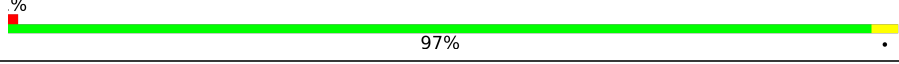

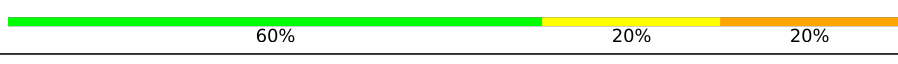
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2808 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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 $\leq 5\%$. 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	250	
1	O	250	
2	B	258	
2	P	258	
3	C	254	
3	Q	254	

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Mol	Chain	Length	Quality of chain
4	D	260	 85% 5% 10%
4	R	260	 85% 5% 10%
5	E	234	 94% 5%
5	S	234	 94% 5%
6	F	288	 81% 16%
6	T	288	 81% 16%
7	G	252	 91% 5%
7	U	252	 91% 5%
8	H	232	 96%
8	V	232	 96%
9	I	205	 98%
9	W	205	 99%
10	J	198	 96%
10	X	198	 95%
11	K	212	 97%
11	Y	212	 97%
12	L	222	 97%
12	Z	222	 97%
13	M	246	 92% 5%
13	a	246	 92% 5%
14	N	196	 97%
14	b	196	 97%
15	c	5	 60% 20% 20%
15	d	5	 60% 20% 20%

2 Entry composition

There are 18 unique types of molecules in this entry. The entry contains 49911 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 Proteasome subunit alpha type-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	250	Total 1915	C 1219	N 315	O 377	S 4	0	0	0
1	O	250	Total 1915	C 1219	N 315	O 377	S 4	0	0	0

- Molecule 2 is a protein called Proteasome subunit alpha type-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	244	Total 1904	C 1201	N 321	O 379	S 3	0	0	0
2	P	244	Total 1904	C 1201	N 321	O 379	S 3	0	0	0

- Molecule 3 is a protein called Proteasome subunit alpha type-4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	C	240	Total 1881	C 1176	N 329	O 372	S 4	0	0	0
3	Q	240	Total 1881	C 1176	N 329	O 372	S 4	0	0	0

- Molecule 4 is a protein called Proteasome subunit alpha type-5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
4	D	235	Total 1813	C 1136	N 304	O 366	S 7	0	0	0
4	R	235	Total 1813	C 1136	N 304	O 366	S 7	0	0	0

- Molecule 5 is a protein called Proteasome subunit alpha type-6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
5	E	231	Total 1773	C 1114	N 307	O 348	S 4	0	0	0
5	S	231	Total 1773	C 1114	N 307	O 348	S 4	0	0	0

- Molecule 6 is a protein called Probable proteasome subunit alpha type-7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
6	F	243	Total 1892	C 1203	N 329	O 356	S 4	0	0	0
6	T	243	Total 1892	C 1203	N 329	O 356	S 4	0	0	0

- Molecule 7 is a protein called Proteasome subunit alpha type-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
7	G	241	Total 1907	C 1214	N 320	O 365	S 8	0	0	0
7	U	241	Total 1907	C 1214	N 320	O 365	S 8	0	0	0

- Molecule 8 is a protein called Proteasome subunit beta type-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
8	H	226	Total 1717	C 1080	N 297	O 333	S 7	0	0	0
8	V	226	Total 1717	C 1080	N 297	O 333	S 7	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
H	116	ASN	HIS	conflict	UNP P25043
V	116	ASN	HIS	conflict	UNP P25043

- Molecule 9 is a protein called Proteasome subunit beta type-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
9	I	204	Total 1581	C 1010	N 258	O 305	S 8	0	0	0
9	W	204	Total 1581	C 1010	N 258	O 305	S 8	0	0	0

- Molecule 10 is a protein called Proteasome subunit beta type-4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	J	195	Total	C	N	O	S	0	0	0
			1561	992	264	299	6			
10	X	195	Total	C	N	O	S	0	0	0
			1561	992	264	299	6			

- Molecule 11 is a protein called Proteasome subunit beta type-5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
11	K	212	Total	C	N	O	S	0	0	0
			1644	1045	280	312	7			
11	Y	212	Total	C	N	O	S	0	0	0
			1644	1045	280	312	7			

- Molecule 12 is a protein called Proteasome subunit beta type-6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
12	L	222	Total	C	N	O	S	0	0	0
			1757	1115	303	335	4			
12	Z	222	Total	C	N	O	S	0	0	0
			1757	1115	303	335	4			

- Molecule 13 is a protein called Proteasome subunit beta type-7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
13	M	233	Total	C	N	O	S	0	0	0
			1824	1154	312	351	7			
13	a	233	Total	C	N	O	S	0	0	0
			1824	1154	312	351	7			

- Molecule 14 is a protein called Proteasome subunit beta type-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
14	N	196	Total	C	N	O	S	0	0	0
			1512	955	250	300	7			
14	b	196	Total	C	N	O	S	0	0	0
			1512	955	250	300	7			

- Molecule 15 is a protein called Ac-PAE-ep.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
15	c	5	Total	C	N	O	0	0	0
			28	18	3	7			
15	d	5	Total	C	N	O	0	0	0
			28	18	3	7			

- Molecule 16 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
16	G	1	Total	Mg	0	0
			1	1		
16	I	2	Total	Mg	0	0
			2	2		
16	K	1	Total	Mg	0	0
			1	1		
16	L	1	Total	Mg	0	0
			1	1		
16	N	1	Total	Mg	0	0
			1	1		
16	Z	1	Total	Mg	0	0
			1	1		

- Molecule 17 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
17	G	1	Total	Cl	0	0
			1	1		
17	U	1	Total	Cl	0	0
			1	1		

- Molecule 18 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
18	A	20	Total	O	0	0
			20	20		
18	B	16	Total	O	0	0
			16	16		
18	C	15	Total	O	0	0
			15	15		
18	D	6	Total	O	0	0
			6	6		
18	E	9	Total	O	0	0
			9	9		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
18	F	12	Total O 12 12	0	0
18	G	29	Total O 29 29	0	0
18	H	27	Total O 27 27	0	0
18	I	13	Total O 13 13	0	0
18	J	21	Total O 21 21	0	0
18	K	20	Total O 20 20	0	0
18	L	30	Total O 30 30	0	0
18	M	23	Total O 23 23	0	0
18	N	17	Total O 17 17	0	0
18	O	11	Total O 11 11	0	0
18	P	8	Total O 8 8	0	0
18	Q	9	Total O 9 9	0	0
18	R	9	Total O 9 9	0	0
18	S	10	Total O 10 10	0	0
18	T	19	Total O 19 19	0	0
18	U	29	Total O 29 29	0	0
18	V	21	Total O 21 21	0	0
18	W	14	Total O 14 14	0	0
18	X	16	Total O 16 16	0	0
18	Y	21	Total O 21 21	0	0
18	Z	18	Total O 18 18	0	0

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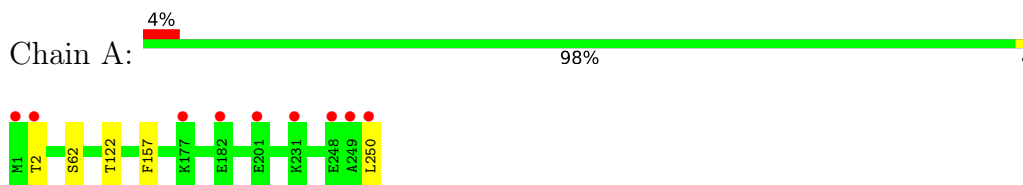
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
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18	b	14	Total 14	O 14	0	0
18	d	1	Total 1	O 1	0	0

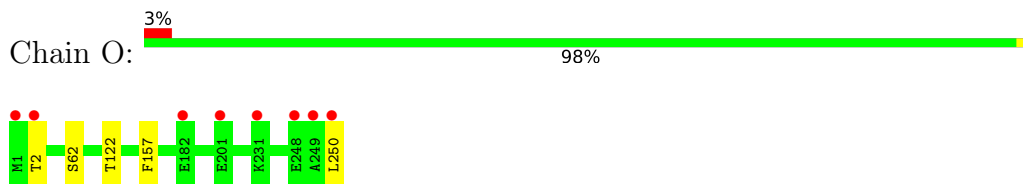
3 Residue-property plots [i](#)

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 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.

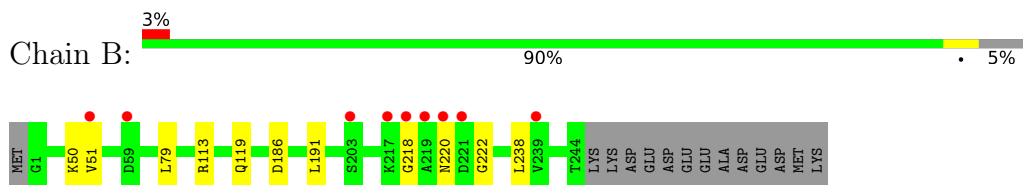
- Molecule 1: Proteasome subunit alpha type-2



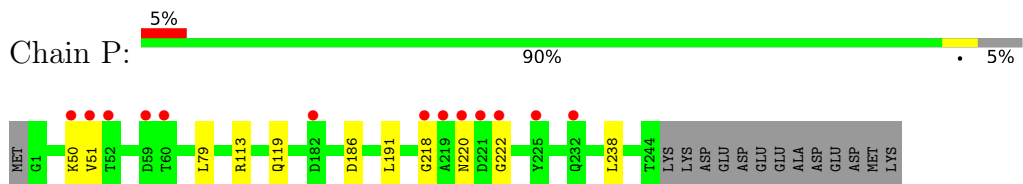
- Molecule 1: Proteasome subunit alpha type-2



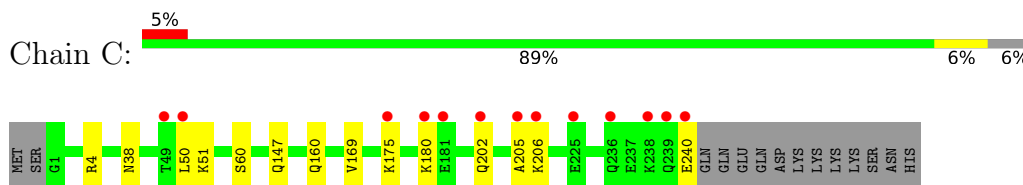
- Molecule 2: Proteasome subunit alpha type-3



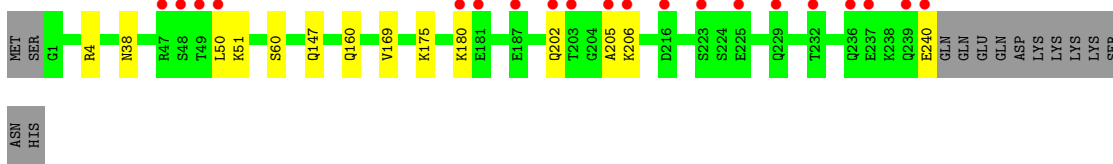
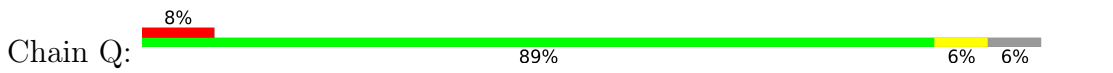
- Molecule 2: Proteasome subunit alpha type-3



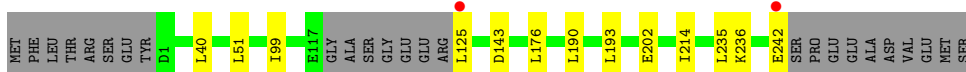
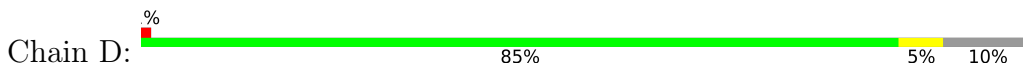
- Molecule 3: Proteasome subunit alpha type-4



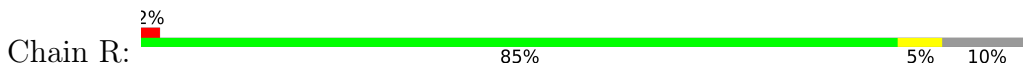
- Molecule 3: Proteasome subunit alpha type-4



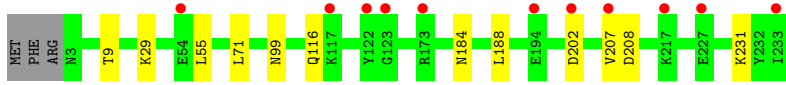
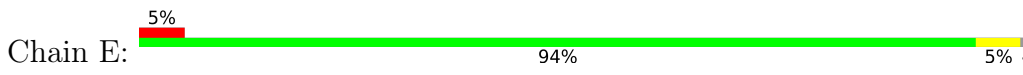
• Molecule 4: Proteasome subunit alpha type-5



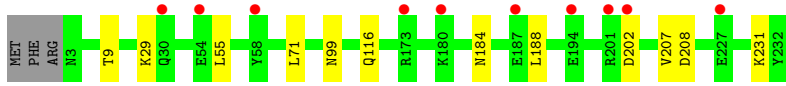
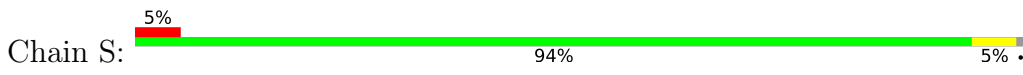
• Molecule 4: Proteasome subunit alpha type-5



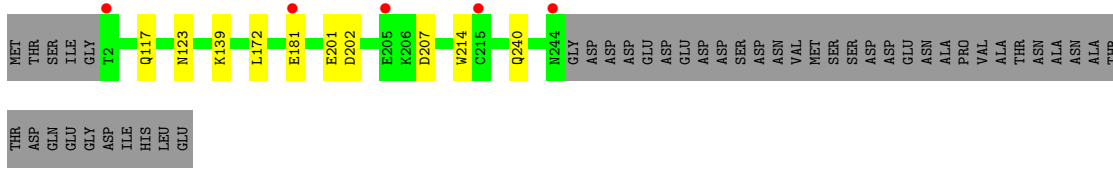
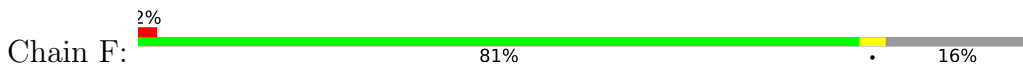
• Molecule 5: Proteasome subunit alpha type-6



• Molecule 5: Proteasome subunit alpha type-6

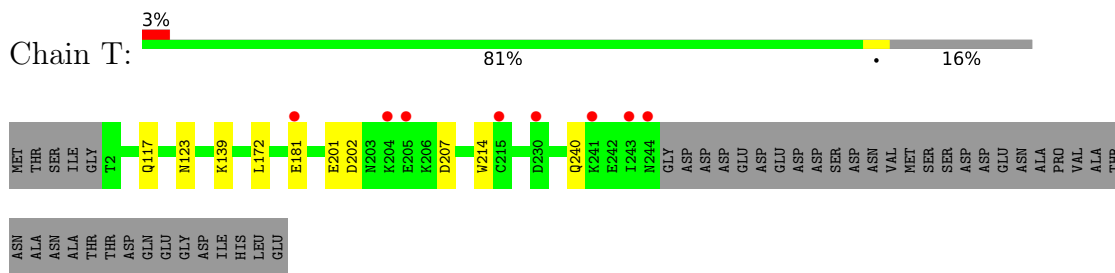


• Molecule 6: Probable proteasome subunit alpha type-7

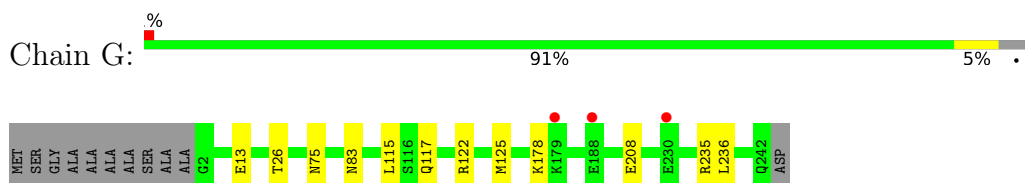


• Molecule 6: Probable proteasome subunit alpha type-7

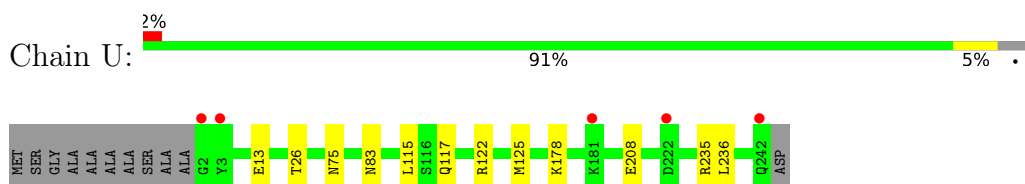




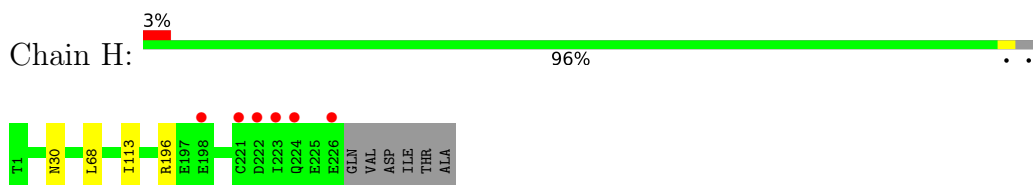
• Molecule 7: Proteasome subunit alpha type-1



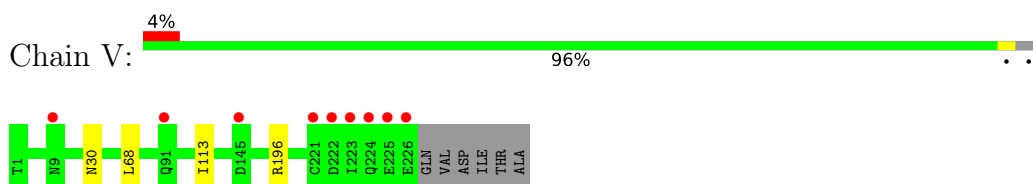
• Molecule 7: Proteasome subunit alpha type-1



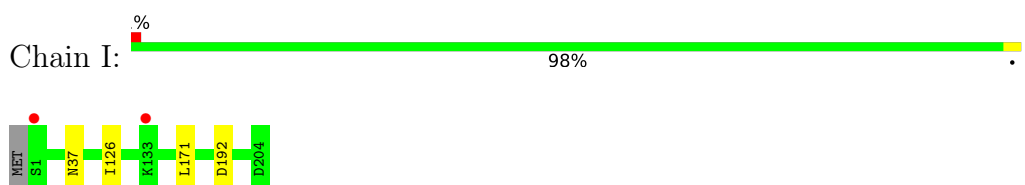
• Molecule 8: Proteasome subunit beta type-2



• Molecule 8: Proteasome subunit beta type-2



• Molecule 9: Proteasome subunit beta type-3

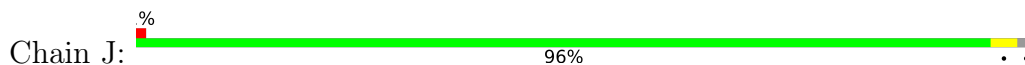


• Molecule 9: Proteasome subunit beta type-3





- Molecule 10: Proteasome subunit beta type-4



- Molecule 10: Proteasome subunit beta type-4



- Molecule 11: Proteasome subunit beta type-5



- Molecule 11: Proteasome subunit beta type-5



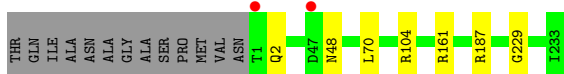
- Molecule 12: Proteasome subunit beta type-6



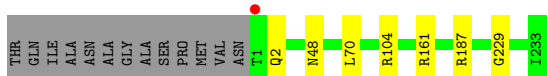
- Molecule 12: Proteasome subunit beta type-6



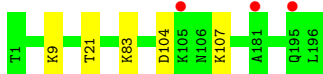
- Molecule 13: Proteasome subunit beta type-7



- Molecule 13: Proteasome subunit beta type-7



- Molecule 14: Proteasome subunit beta type-1



- Molecule 14: Proteasome subunit beta type-1



- Molecule 15: Ac-PAE-ep



- Molecule 15: Ac-PAE-ep



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	137.50Å 301.36Å 145.84Å 90.00° 113.25° 90.00°	Depositor
Resolution (Å)	15.00 – 2.70 15.00 – 2.70	Depositor EDS
% Data completeness (in resolution range)	96.9 (15.00-2.70) 96.9 (15.00-2.70)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.60 (at 2.69Å)	Xtrriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.202 , 0.220 0.205 , 0.223	Depositor DCC
R_{free} test set	14322 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	58.1	Xtrriage
Anisotropy	0.072	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 39.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	49911	wwPDB-VP
Average B, all atoms (Å ²)	65.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.28% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GAU, MG, POL, ACE, CL

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	0.29	0/1952	0.47	0/2642
1	O	0.29	0/1952	0.47	0/2642
2	B	0.29	0/1934	0.50	0/2618
2	P	0.29	0/1934	0.51	0/2618
3	C	0.29	0/1910	0.52	0/2586
3	Q	0.29	0/1910	0.51	0/2586
4	D	0.28	0/1837	0.49	0/2475
4	R	0.28	0/1837	0.49	0/2475
5	E	0.28	0/1800	0.48	0/2433
5	S	0.28	0/1800	0.48	0/2433
6	F	0.29	0/1932	0.46	0/2609
6	T	0.28	0/1932	0.46	0/2609
7	G	0.29	0/1945	0.47	0/2634
7	U	0.28	0/1945	0.47	0/2634
8	H	0.30	0/1747	0.49	0/2369
8	V	0.28	0/1747	0.48	0/2369
9	I	0.30	0/1611	0.49	0/2174
9	W	0.31	0/1611	0.49	0/2174
10	J	0.30	0/1589	0.50	0/2142
10	X	0.28	0/1589	0.49	0/2142
11	K	0.29	0/1681	0.51	1/2274 (0.0%)
11	Y	0.27	0/1681	0.51	1/2274 (0.0%)
12	L	0.29	0/1795	0.50	0/2420
12	Z	0.29	0/1795	0.50	0/2420
13	M	0.29	0/1855	0.51	0/2514
13	a	0.29	0/1855	0.51	0/2514
14	N	0.35	0/1541	0.51	0/2087
14	b	0.34	0/1541	0.50	0/2087
15	c	2.69	1/13 (7.7%)	1.44	0/18
15	d	2.62	1/13 (7.7%)	1.34	0/18
All	All	0.30	2/50284 (0.0%)	0.49	2/67990 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
15	d	2	PRO	CA-C	-7.67	1.37	1.52
15	c	2	PRO	CA-C	-7.31	1.38	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	K	4	LEU	CA-CB-CG	5.18	127.20	115.30
11	Y	4	LEU	CA-CB-CG	5.17	127.18	115.30

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 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	248/250 (99%)	238 (96%)	9 (4%)	1 (0%)	34	60
1	O	248/250 (99%)	238 (96%)	9 (4%)	1 (0%)	34	60
2	B	242/258 (94%)	235 (97%)	3 (1%)	4 (2%)	9	23
2	P	242/258 (94%)	234 (97%)	4 (2%)	4 (2%)	9	23
3	C	238/254 (94%)	233 (98%)	3 (1%)	2 (1%)	19	43
3	Q	238/254 (94%)	233 (98%)	3 (1%)	2 (1%)	19	43
4	D	231/260 (89%)	228 (99%)	3 (1%)	0	100	100
4	R	231/260 (89%)	228 (99%)	3 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
5	E	229/234 (98%)	222 (97%)	7 (3%)	0	100	100
5	S	229/234 (98%)	221 (96%)	8 (4%)	0	100	100
6	F	241/288 (84%)	238 (99%)	3 (1%)	0	100	100
6	T	241/288 (84%)	238 (99%)	3 (1%)	0	100	100
7	G	239/252 (95%)	238 (100%)	1 (0%)	0	100	100
7	U	239/252 (95%)	238 (100%)	1 (0%)	0	100	100
8	H	224/232 (97%)	218 (97%)	6 (3%)	0	100	100
8	V	224/232 (97%)	218 (97%)	6 (3%)	0	100	100
9	I	202/205 (98%)	196 (97%)	6 (3%)	0	100	100
9	W	202/205 (98%)	196 (97%)	6 (3%)	0	100	100
10	J	193/198 (98%)	189 (98%)	3 (2%)	1 (0%)	29	54
10	X	193/198 (98%)	189 (98%)	3 (2%)	1 (0%)	29	54
11	K	210/212 (99%)	206 (98%)	4 (2%)	0	100	100
11	Y	210/212 (99%)	206 (98%)	4 (2%)	0	100	100
12	L	220/222 (99%)	215 (98%)	5 (2%)	0	100	100
12	Z	220/222 (99%)	215 (98%)	5 (2%)	0	100	100
13	M	231/246 (94%)	224 (97%)	6 (3%)	1 (0%)	34	60
13	a	231/246 (94%)	224 (97%)	6 (3%)	1 (0%)	34	60
14	N	194/196 (99%)	189 (97%)	5 (3%)	0	100	100
14	b	194/196 (99%)	189 (97%)	5 (3%)	0	100	100
15	c	2/5 (40%)	2 (100%)	0	0	100	100
15	d	2/5 (40%)	2 (100%)	0	0	100	100
All	All	6288/6624 (95%)	6140 (98%)	130 (2%)	18 (0%)	41	66

5 of 18 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	51	VAL
2	B	222	GLY
3	C	202	GLN
2	P	51	VAL
2	P	222	GLY

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	209/209 (100%)	205 (98%)	4 (2%)	57 82
1	O	209/209 (100%)	205 (98%)	4 (2%)	57 82
2	B	203/216 (94%)	196 (97%)	7 (3%)	37 66
2	P	203/216 (94%)	196 (97%)	7 (3%)	37 66
3	C	212/226 (94%)	200 (94%)	12 (6%)	20 44
3	Q	212/226 (94%)	200 (94%)	12 (6%)	20 44
4	D	194/215 (90%)	181 (93%)	13 (7%)	16 37
4	R	194/215 (90%)	181 (93%)	13 (7%)	16 37
5	E	190/193 (98%)	178 (94%)	12 (6%)	18 40
5	S	190/193 (98%)	178 (94%)	12 (6%)	18 40
6	F	201/239 (84%)	191 (95%)	10 (5%)	24 51
6	T	201/239 (84%)	191 (95%)	10 (5%)	24 51
7	G	206/210 (98%)	194 (94%)	12 (6%)	20 43
7	U	206/210 (98%)	194 (94%)	12 (6%)	20 43
8	H	185/190 (97%)	181 (98%)	4 (2%)	52 79
8	V	185/190 (97%)	181 (98%)	4 (2%)	52 79
9	I	172/173 (99%)	168 (98%)	4 (2%)	50 78
9	W	172/173 (99%)	170 (99%)	2 (1%)	71 88
10	J	173/175 (99%)	169 (98%)	4 (2%)	50 78
10	X	173/175 (99%)	168 (97%)	5 (3%)	42 71
11	K	169/169 (100%)	162 (96%)	7 (4%)	30 59
11	Y	169/169 (100%)	162 (96%)	7 (4%)	30 59
12	L	185/185 (100%)	178 (96%)	7 (4%)	33 62
12	Z	185/185 (100%)	178 (96%)	7 (4%)	33 62
13	M	199/208 (96%)	193 (97%)	6 (3%)	41 70
13	a	199/208 (96%)	193 (97%)	6 (3%)	41 70

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
14	N	162/162 (100%)	157 (97%)	5 (3%)	40	69
14	b	162/162 (100%)	157 (97%)	5 (3%)	40	69
15	c	1/1 (100%)	0	1 (100%)	0	0
15	d	1/1 (100%)	0	1 (100%)	0	0
All	All	5322/5542 (96%)	5107 (96%)	215 (4%)	31	60

5 of 215 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	P	186	ASP
5	S	9	THR
12	Z	132	GLU
3	Q	38	ASN
4	R	40	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 116 such sidechains are listed below:

Mol	Chain	Res	Type
13	M	102	GLN
12	Z	70	ASN
3	Q	160	GLN
12	Z	49	ASN
8	V	114	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
15	GAU	c	4	15,14	8,8,8	1.50	2 (25%)	8,9,9	1.51	2 (25%)
15	GAU	d	4	15,14	8,8,8	1.45	2 (25%)	8,9,9	1.51	2 (25%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
15	GAU	c	4	15,14	-	0/7/7/7	-
15	GAU	d	4	15,14	-	2/7/7/7	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
15	c	4	GAU	C-CA	2.94	1.57	1.52
15	d	4	GAU	C-CA	2.65	1.56	1.52
15	c	4	GAU	OE2-CD	-2.46	1.22	1.30
15	d	4	GAU	OE2-CD	-2.44	1.22	1.30

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
15	d	4	GAU	CB-CA-N	-2.79	100.94	109.03
15	c	4	GAU	CB-CA-N	-2.72	101.12	109.03
15	c	4	GAU	CG-CB-CA	-2.31	107.90	112.74
15	d	4	GAU	CG-CB-CA	-2.10	108.33	112.74

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
15	d	4	GAU	O-C-CA-CB
15	d	4	GAU	O-C-CA-N

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 9 are 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](#)

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	250/250 (100%)	-0.31	9 (3%) 42 42	40, 53, 88, 130	0
1	O	250/250 (100%)	-0.22	8 (3%) 47 48	46, 63, 107, 134	0
2	B	244/258 (94%)	-0.18	9 (3%) 41 41	41, 59, 101, 151	0
2	P	244/258 (94%)	-0.04	13 (5%) 26 25	49, 66, 110, 158	0
3	C	240/254 (94%)	-0.08	13 (5%) 25 24	42, 64, 126, 151	0
3	Q	240/254 (94%)	0.15	20 (8%) 11 9	50, 77, 151, 174	0
4	D	235/260 (90%)	-0.25	2 (0%) 84 85	49, 67, 98, 139	0
4	R	235/260 (90%)	-0.13	6 (2%) 56 57	51, 71, 110, 139	0
5	E	231/234 (98%)	-0.06	11 (4%) 30 28	50, 70, 109, 157	0
5	S	231/234 (98%)	0.07	11 (4%) 30 28	53, 77, 120, 166	0
6	F	243/288 (84%)	-0.32	5 (2%) 63 65	44, 61, 109, 137	0
6	T	243/288 (84%)	-0.11	8 (3%) 46 46	48, 72, 126, 156	0
7	G	241/252 (95%)	-0.41	3 (1%) 79 80	36, 55, 87, 137	0
7	U	241/252 (95%)	-0.29	5 (2%) 63 65	44, 59, 91, 133	0
8	H	226/232 (97%)	-0.27	6 (2%) 54 55	38, 53, 86, 149	0
8	V	226/232 (97%)	-0.22	9 (3%) 38 37	44, 56, 89, 170	0
9	I	204/205 (99%)	-0.59	2 (0%) 82 83	39, 51, 79, 106	0
9	W	204/205 (99%)	-0.46	3 (1%) 73 76	43, 55, 84, 109	0
10	J	195/198 (98%)	-0.46	2 (1%) 82 83	41, 55, 78, 133	0
10	X	195/198 (98%)	-0.41	2 (1%) 82 83	42, 57, 80, 143	0
11	K	212/212 (100%)	-0.43	2 (0%) 84 85	39, 57, 79, 100	0
11	Y	212/212 (100%)	-0.42	0 100 100	43, 56, 79, 99	0
12	L	222/222 (100%)	-0.42	2 (0%) 84 85	41, 56, 90, 123	0
12	Z	222/222 (100%)	-0.42	4 (1%) 68 70	36, 55, 87, 119	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
13	M	233/246 (94%)	-0.49	2 (0%) 84 85	39, 54, 75, 96	0
13	a	233/246 (94%)	-0.44	1 (0%) 92 93	38, 54, 74, 93	0
14	N	196/196 (100%)	-0.52	3 (1%) 73 76	37, 49, 76, 104	0
14	b	196/196 (100%)	-0.50	2 (1%) 82 83	39, 51, 76, 111	0
15	c	2/5 (40%)	-0.64	0 100 100	60, 60, 60, 63	0
15	d	2/5 (40%)	-1.01	0 100 100	54, 54, 54, 59	0
All	All	6348/6624 (95%)	-0.29	163 (2%) 56 57	36, 60, 104, 174	0

The worst 5 of 163 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	P	218	GLY	7.0
10	X	1	MET	6.5
3	Q	49	THR	6.5
3	Q	50	LEU	6.4
3	Q	206	LYS	6.3

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
15	GAU	d	4	9/9	0.96	0.15	53,61,64,64	0
15	GAU	c	4	9/9	0.97	0.12	57,66,71,72	0

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
16	MG	N	201	1/1	0.93	0.12	53,53,53,53	0
16	MG	G	301	1/1	0.95	0.10	51,51,51,51	0
17	CL	U	301	1/1	0.95	0.27	30,30,30,30	0
16	MG	Z	301	1/1	0.96	0.30	60,60,60,60	0
16	MG	I	301	1/1	0.96	0.38	67,67,67,67	0
17	CL	G	302	1/1	0.97	0.19	30,30,30,30	0
16	MG	L	301	1/1	0.98	0.05	67,67,67,67	0
16	MG	K	301	1/1	0.99	0.07	49,49,49,49	0
16	MG	I	302	1/1	0.99	0.06	53,53,53,53	0

6.5 Other polymers [i](#)

There are no such residues in this entry.