



# Full wwPDB X-ray Structure Validation Report ⓘ

Dec 12, 2016 – 06:34 PM EST

PDB ID : 5D0S  
Title : Yeast 20S proteasome beta5-D166N mutant in complex with Carfilzomib  
Authors : Huber, E.M.; Groll, M.  
Deposited on : 2015-08-03  
Resolution : 2.50 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](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.7.1 (RC1), CSD as537be (2016)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20028442  
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) : rb-20028442

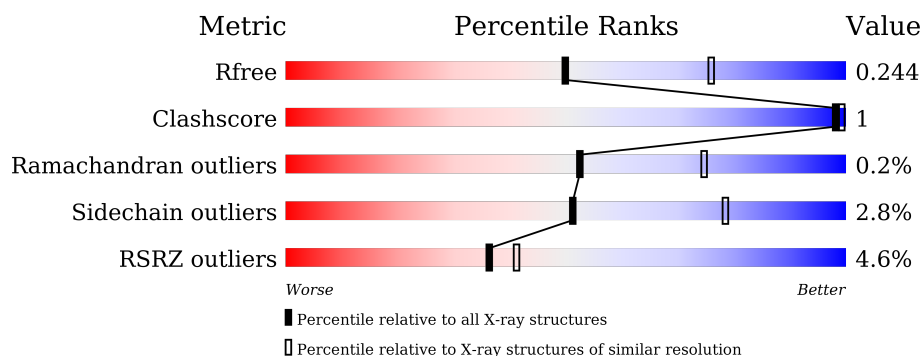
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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  $\geq 3$ , 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	250	<div> <div>5%</div> <div>98%</div> <div>.</div> </div>
1	O	250	<div> <div>6%</div> <div>98%</div> <div>.</div> </div>
2	B	258	<div> <div>8%</div> <div>90%</div> <div>.. 5%</div> </div>
2	P	258	<div> <div>8%</div> <div>91%</div> <div>.. 5%</div> </div>
3	C	254	<div> <div>11%</div> <div>90%</div> <div>.. 6%</div> </div>
3	Q	254	<div> <div>15%</div> <div>90%</div> <div>.. 6%</div> </div>

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Mol	Chain	Length	Quality of chain
4	D	260	
4	R	260	
5	E	234	
5	S	234	
6	F	288	
6	T	288	
7	G	252	
7	U	252	
8	H	232	
8	V	232	
9	I	205	
9	W	205	
10	J	198	
10	X	198	
11	K	212	
11	Y	212	
12	L	222	
12	Z	222	
13	M	246	
13	a	246	
14	N	196	
14	b	196	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
15	MG	Z	301	-	-	-	X
17	3BV	N	201	-	-	-	X
17	3BV	b	201	-	-	-	X
18	MES	H	302	-	-	-	X
18	MES	K	303	-	-	-	X
18	MES	V	303	-	-	-	X
18	MES	Y	303	-	-	-	X

## 2 Entry composition

There are 19 unique types of molecules in this entry. The entry contains 50766 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
1	A	250	Total	C	N	O	S	0	0	0
			1915	1219	315	377	4			
1	O	250	Total	C	N	O	S	0	0	0
			1915	1219	315	377	4			

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

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

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

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

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

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

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

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

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

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

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	G	241	Total	C	N	O	S	0	0	0
			1907	1214	320	365	8			
7	U	241	Total	C	N	O	S	0	0	0
			1906	1214	320	364	8			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	H	222	Total	C	N	O	S	0	0	0
			1684	1061	293	323	7			
8	V	222	Total	C	N	O	S	0	0	0
			1684	1061	293	323	7			

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

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

- 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			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
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	281	311	7			
11	Y	212	Total	C	N	O	S	0	0	0
			1644	1045	281	311	7			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
K	168	ASN	ASP	engineered mutation	UNP P30656
Y	168	ASN	ASP	engineered mutation	UNP P30656

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

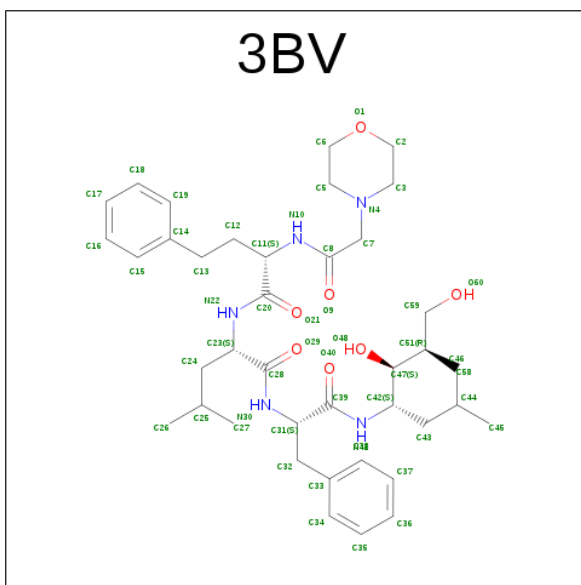
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
15	G	1	Total Mg 1 1	0	0
15	K	1	Total Mg 1 1	0	0
15	I	1	Total Mg 1 1	0	0
15	V	1	Total Mg 1 1	0	0
15	Z	1	Total Mg 1 1	0	0
15	N	1	Total Mg 1 1	0	0
15	Y	1	Total Mg 1 1	0	0

- Molecule 16 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
16	G	1	Total Cl 1 1	0	0
16	b	1	Total Cl 1 1	0	0
16	N	1	Total Cl 1 1	0	0
16	U	1	Total Cl 1 1	0	0

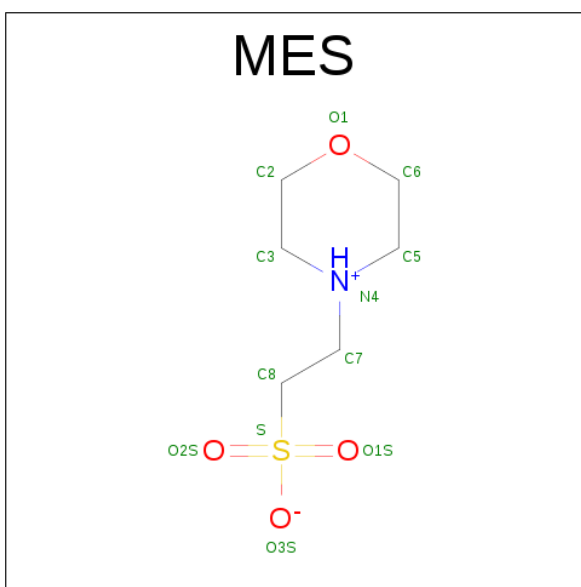
- Molecule 17 is N-[(2S)-2-[(morpholin-4-ylacetyl)amino]-4-phenylbutanoyl]-L-leucyl-N-[(2R,3S,4S)-1,3-dihydroxy-2,6-dimethylheptan-4-yl]-L-phenylalaninamide (three-letter code: 3BV) (formula: C<sub>40</sub>H<sub>61</sub>N<sub>5</sub>O<sub>7</sub>).





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
17	H	1	Total 52	C 40	N 5	O 7	0	0
17	K	1	Total 52	C 40	N 5	O 7	0	0
17	N	1	Total 52	C 40	N 5	O 7	0	0
17	V	1	Total 52	C 40	N 5	O 7	0	0
17	Y	1	Total 52	C 40	N 5	O 7	0	0
17	b	1	Total 52	C 40	N 5	O 7	0	0

- Molecule 18 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula:  $\text{C}_6\text{H}_{13}\text{NO}_4\text{S}$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
18	H	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
18	K	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
18	V	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
18	Y	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 19 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
19	A	52	Total	O	0	0
			52	52		
19	B	31	Total	O	0	0
			31	31		
19	C	38	Total	O	0	0
			38	38		
19	D	31	Total	O	0	0
			31	31		
19	E	27	Total	O	0	0
			27	27		
19	F	38	Total	O	0	0
			38	38		
19	G	47	Total	O	0	0
			47	47		
19	H	57	Total	O	0	0
			57	57		

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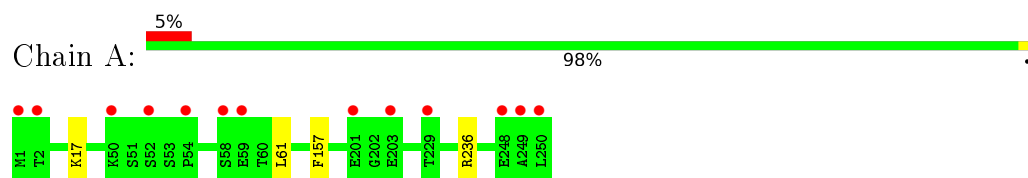
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
19	I	49	Total 49	O 49	0	0
19	J	52	Total 52	O 52	0	0
19	K	54	Total 54	O 54	0	0
19	L	54	Total 54	O 54	0	0
19	M	48	Total 48	O 48	0	0
19	N	43	Total 43	O 43	0	0
19	O	27	Total 27	O 27	0	0
19	P	29	Total 29	O 29	0	0
19	Q	15	Total 15	O 15	0	0
19	R	18	Total 18	O 18	0	0
19	S	11	Total 11	O 11	0	0
19	T	28	Total 28	O 28	0	0
19	U	44	Total 44	O 44	0	0
19	V	42	Total 42	O 42	0	0
19	W	35	Total 35	O 35	0	0
19	X	47	Total 47	O 47	0	0
19	Y	35	Total 35	O 35	0	0
19	Z	47	Total 47	O 47	0	0
19	a	64	Total 64	O 64	0	0
19	b	37	Total 37	O 37	0	0

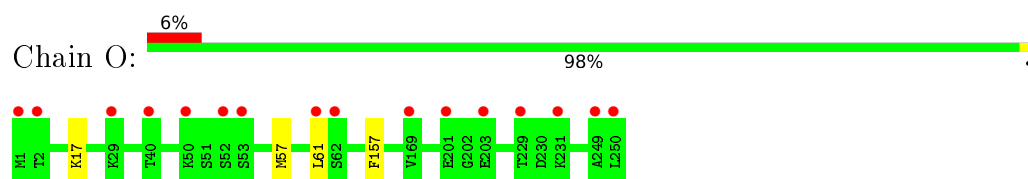
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

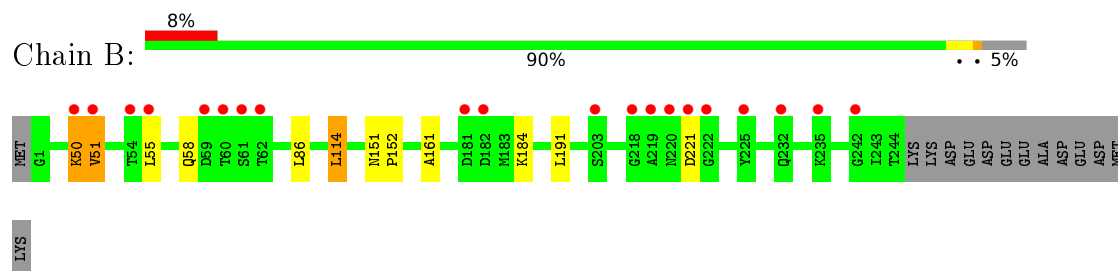
- Molecule 1: Proteasome subunit alpha type-2



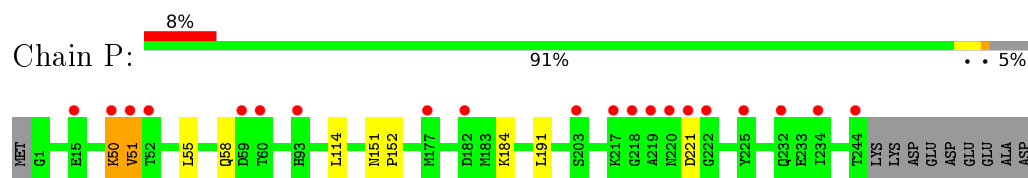
- Molecule 1: Proteasome subunit alpha type-2



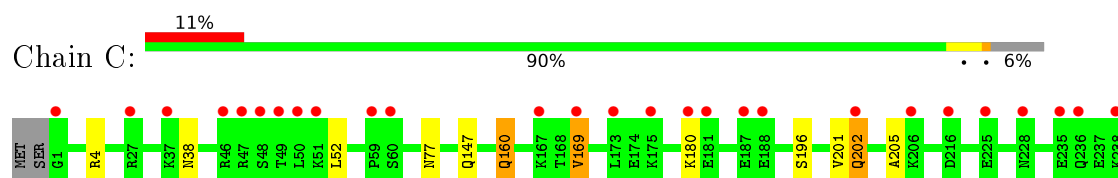
- Molecule 2: Proteasome subunit alpha type-3

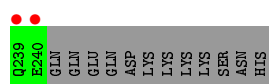


- Molecule 2: Proteasome subunit alpha type-3

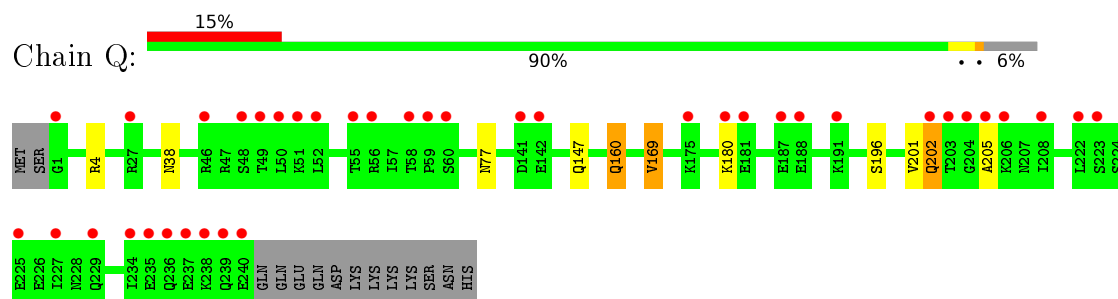


- Molecule 3: Proteasome subunit alpha type-4

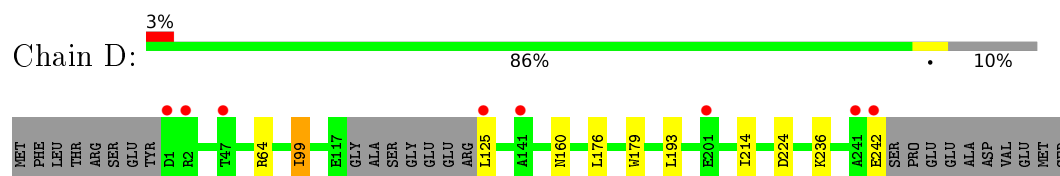




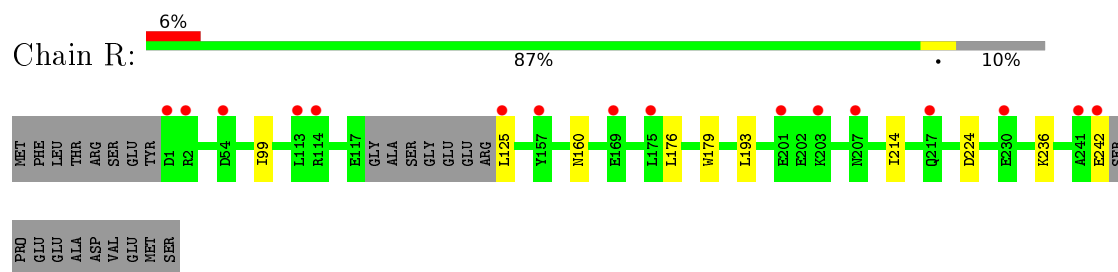
- Molecule 3: Proteasome subunit alpha type-4



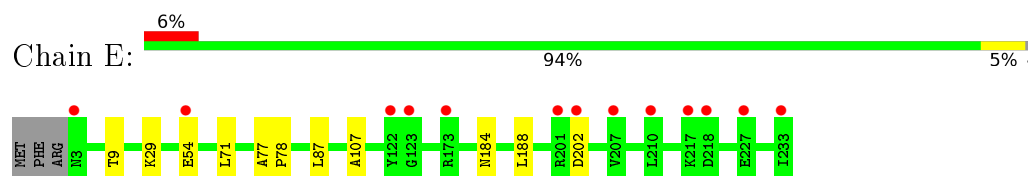
- Molecule 4: Proteasome subunit alpha type-5



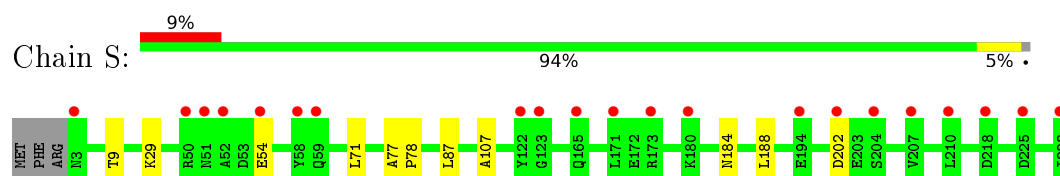
- Molecule 4: Proteasome subunit alpha type-5



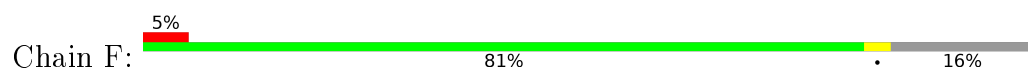
- Molecule 5: Proteasome subunit alpha type-6

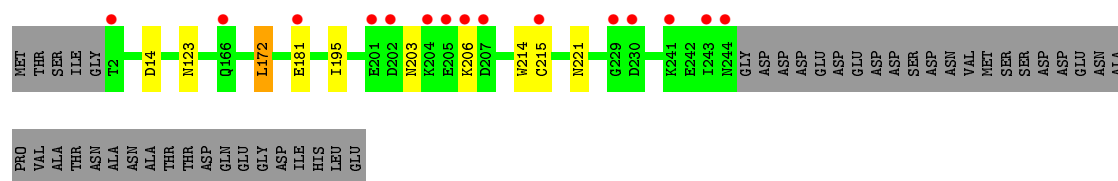


- Molecule 5: Proteasome subunit alpha type-6

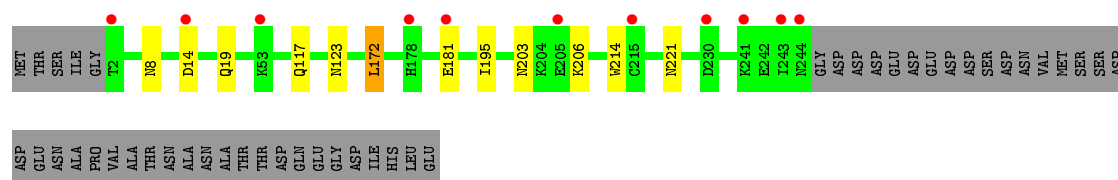
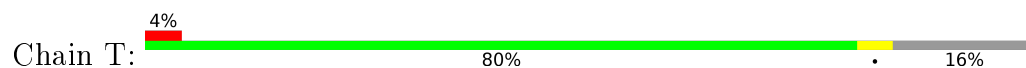


- Molecule 6: Probable proteasome subunit alpha type-7

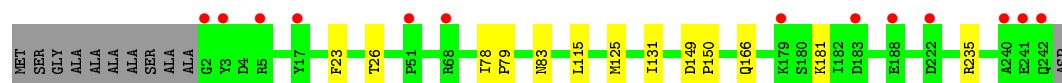
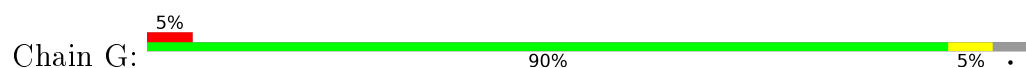




- Molecule 6: Probable proteasome subunit alpha type-7



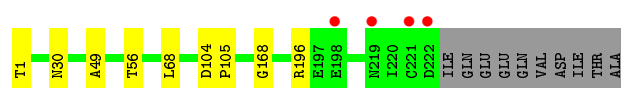
- Molecule 7: Proteasome subunit alpha type-1



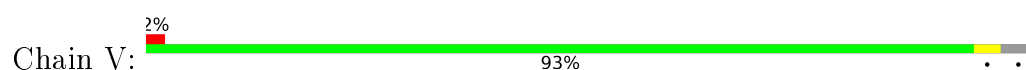
- Molecule 7: Proteasome subunit alpha type-1



- Molecule 8: Proteasome subunit beta type-2



- Molecule 8: Proteasome subunit beta type-2



- Molecule 9: Proteasome subunit beta type-3





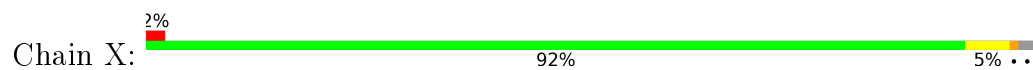
- Molecule 9: Proteasome subunit beta type-3



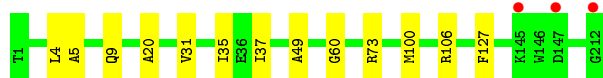
- Molecule 10: Proteasome subunit beta type-4



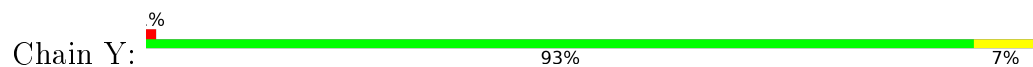
- Molecule 10: Proteasome subunit beta type-4



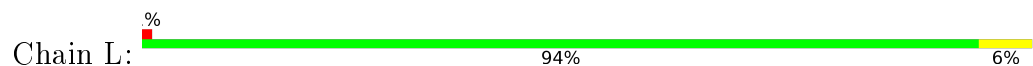
- Molecule 11: Proteasome subunit beta type-5



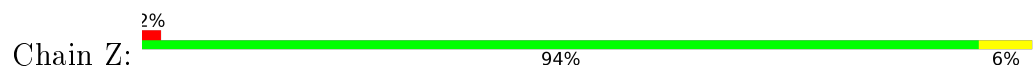
- Molecule 11: Proteasome subunit beta type-5



- Molecule 12: Proteasome subunit beta type-6

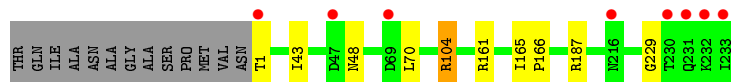


- Molecule 12: Proteasome subunit beta type-6

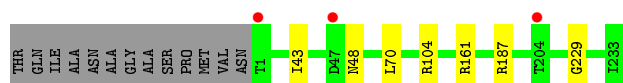
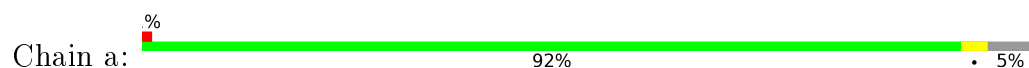




- Molecule 13: Proteasome subunit beta type-7



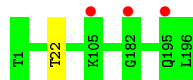
- Molecule 13: Proteasome subunit beta type-7



- Molecule 14: Proteasome subunit beta type-1



- Molecule 14: Proteasome subunit beta type-1





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	135.02Å 300.64Å 144.31Å 90.00° 113.02° 90.00°	Depositor
Resolution (Å)	15.00 – 2.50 15.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	96.4 (15.00-2.50) 96.4 (15.00-2.50)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.27 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, $R_{free}$	0.216 , 0.237 0.224 , 0.244	Depositor DCC
$R_{free}$ test set	17448 reflections (5.26%)	DCC
Wilson B-factor (Å <sup>2</sup> )	43.2	Xtriage
Anisotropy	0.088	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 37.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	50766	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.35% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG, 3BV, MES, 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	0.27	0/1952	0.46	0/2642
1	O	0.27	0/1952	0.46	0/2642
2	B	0.27	0/1934	0.48	0/2618
2	P	0.26	0/1934	0.48	0/2618
3	C	0.27	0/1910	0.50	0/2586
3	Q	0.27	0/1910	0.49	0/2586
4	D	0.26	0/1837	0.46	0/2475
4	R	0.26	0/1837	0.46	0/2475
5	E	0.26	0/1800	0.46	0/2433
5	S	0.26	0/1800	0.46	0/2433
6	F	0.27	0/1932	0.45	0/2609
6	T	0.27	0/1932	0.45	0/2609
7	G	0.27	0/1945	0.47	0/2634
7	U	0.27	0/1944	0.46	0/2632
8	H	0.28	0/1715	0.47	1/2326 (0.0%)
8	V	0.28	0/1715	0.47	0/2326
9	I	0.27	0/1611	0.48	0/2174
9	W	0.27	0/1611	0.48	0/2174
10	J	0.26	0/1589	0.48	0/2142
10	X	0.26	0/1589	0.48	0/2142
11	K	0.33	0/1681	0.50	0/2274
11	Y	0.30	0/1681	0.49	0/2274
12	L	0.27	0/1795	0.47	0/2420
12	Z	0.27	0/1795	0.47	0/2420
13	M	0.27	0/1855	0.50	0/2514
13	a	0.27	0/1855	0.51	0/2514
14	N	0.27	0/1541	0.47	0/2087
14	b	0.27	0/1541	0.48	0/2087
All	All	0.27	0/50193	0.47	1/67866 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	H	1	THR	N-CA-C	5.50	125.85	111.00

There are no chirality outliers.

There are no planarity outliers.

## 5.2 Too-close contacts ⓘ

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	1915	0	1929	1	0
1	O	1915	0	1929	1	0
2	B	1904	0	1904	4	0
2	P	1904	0	1904	2	0
3	C	1881	0	1895	5	0
3	Q	1881	0	1895	4	0
4	D	1813	0	1797	3	0
4	R	1813	0	1797	1	0
5	E	1773	0	1775	2	0
5	S	1773	0	1775	2	0
6	F	1892	0	1883	2	0
6	T	1892	0	1883	3	0
7	G	1907	0	1901	4	0
7	U	1906	0	1901	4	0
8	H	1684	0	1685	3	0
8	V	1684	0	1685	2	0
9	I	1581	0	1574	4	0
9	W	1581	0	1574	4	0
10	J	1561	0	1569	6	0
10	X	1561	0	1569	5	0
11	K	1644	0	1594	7	0
11	Y	1644	0	1594	8	0
12	L	1757	0	1711	8	0
12	Z	1757	0	1711	5	0
13	M	1824	0	1832	3	0
13	a	1824	0	1832	0	0
14	N	1512	0	1478	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
14	b	1512	0	1478	0	0
15	G	1	0	0	0	0
15	I	1	0	0	0	0
15	K	1	0	0	0	0
15	N	1	0	0	0	0
15	V	1	0	0	0	0
15	Y	1	0	0	0	0
15	Z	1	0	0	0	0
16	G	1	0	0	0	0
16	N	1	0	0	0	0
16	U	1	0	0	0	0
16	b	1	0	0	0	0
17	H	52	0	59	2	0
17	K	52	0	58	4	0
17	N	52	0	58	1	0
17	V	52	0	58	4	0
17	Y	52	0	58	2	0
17	b	52	0	58	0	0
18	H	12	0	13	0	0
18	K	12	0	13	1	0
18	V	12	0	13	1	0
18	Y	12	0	13	0	0
19	A	52	0	0	1	0
19	B	31	0	0	0	0
19	C	38	0	0	0	0
19	D	31	0	0	2	0
19	E	27	0	0	0	0
19	F	38	0	0	1	0
19	G	47	0	0	0	0
19	H	57	0	0	0	0
19	I	49	0	0	0	0
19	J	52	0	0	1	0
19	K	54	0	0	0	0
19	L	54	0	0	2	0
19	M	48	0	0	0	0
19	N	43	0	0	1	0
19	O	27	0	0	0	0
19	P	29	0	0	0	0
19	Q	15	0	0	0	0
19	R	18	0	0	0	0
19	S	11	0	0	0	0
19	T	28	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
19	U	44	0	0	1	0
19	V	42	0	0	0	0
19	W	35	0	0	0	0
19	X	47	0	0	1	0
19	Y	35	0	0	0	0
19	Z	47	0	0	0	0
19	a	64	0	0	0	0
19	b	37	0	0	0	0
All	All	50766	0	49455	95	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 1.

All (95) 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)
11:Y:5:ALA:HB3	11:Y:100:MET:HE2	1.68	0.76
11:K:5:ALA:HB3	11:K:100:MET:HE2	1.69	0.75
6:T:19:GLN:NE2	19:T:301:HOH:O	2.23	0.72
14:N:35:THR:HG22	19:N:305:HOH:O	1.90	0.72
17:K:301:3BV:H16	12:L:108:HIS:NE2	2.10	0.67
12:Z:13:LEU:HD13	12:Z:150:LEU:HD21	1.77	0.66
12:L:13:LEU:HD13	12:L:150:LEU:HD21	1.78	0.65
17:N:201:3BV:H1	17:N:201:3BV:O9	1.98	0.63
17:K:301:3BV:H60	18:K:303:MES:O1S	2.05	0.56
6:F:215:CYS:HB3	19:F:304:HOH:O	2.06	0.56
8:H:49:ALA:HA	17:H:301:3BV:H50	1.88	0.56
10:J:67:TYR:CE1	10:J:75:LEU:HD13	2.44	0.52
10:X:67:TYR:CE1	10:X:75:LEU:HD13	2.44	0.52
11:K:5:ALA:HB3	11:K:100:MET:CE	2.40	0.52
17:K:301:3BV:H16	12:L:108:HIS:CE1	2.44	0.52
3:C:201:VAL:O	3:C:202:GLN:CB	2.58	0.51
10:J:16:ALA:HB2	10:J:161:LEU:HD21	1.91	0.51
3:Q:201:VAL:O	3:Q:202:GLN:CB	2.59	0.51
10:X:16:ALA:HB2	10:X:161:LEU:HD21	1.91	0.51
11:Y:170:TYR:O	17:Y:301:3BV:H57	2.10	0.51
11:Y:5:ALA:HB3	11:Y:100:MET:CE	2.39	0.51
11:K:49:ALA:HA	17:K:301:3BV:H50	1.92	0.51
3:C:169:VAL:HG23	3:C:196:SER:HB2	1.94	0.50
14:N:152:VAL:HA	14:N:175:MET:HE1	1.92	0.50
11:Y:100:MET:HE3	11:Y:127:PHE:HB2	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:V:301:3BV:O60	18:V:303:MES:O1S	2.29	0.50
7:G:23:PHE:O	7:G:26:THR:HB	2.12	0.50
3:Q:169:VAL:HG23	3:Q:196:SER:HB2	1.94	0.50
4:D:99:ILE:HG23	19:D:321:HOH:O	2.12	0.49
11:K:100:MET:HE3	11:K:127:PHE:HB2	1.93	0.49
7:U:23:PHE:O	7:U:26:THR:HB	2.12	0.49
11:Y:49:ALA:HA	17:Y:301:3BV:H50	1.95	0.48
12:L:2:PHE:N	19:L:302:HOH:O	2.47	0.48
5:S:87:LEU:HD21	5:S:107:ALA:HB1	1.96	0.47
1:A:236:ARG:NH2	19:A:302:HOH:O	2.48	0.47
5:E:87:LEU:HD21	5:E:107:ALA:HB1	1.97	0.46
8:H:104:ASP:HB2	8:H:105:PRO:HD2	1.96	0.46
11:K:37:ILE:HG23	11:K:60:GLY:HA2	1.98	0.46
17:V:301:3BV:O9	17:V:301:3BV:H16	2.15	0.45
6:T:172:LEU:CD1	6:T:195:ILE:HD13	2.46	0.45
6:F:172:LEU:CD1	6:F:195:ILE:HD13	2.46	0.45
10:J:36:ARG:NH1	10:J:58:GLU:OE2	2.48	0.45
9:I:10:ILE:HG21	9:I:141:ALA:HB3	1.99	0.45
10:X:36:ARG:NH1	10:X:58:GLU:OE2	2.50	0.45
19:L:302:HOH:O	13:M:1:THR:HG23	2.16	0.45
3:Q:160:GLN:HA	3:Q:160:GLN:HE21	1.81	0.45
4:R:160:ASN:HB3	4:R:179:TRP:CE2	2.52	0.44
8:V:104:ASP:HB2	8:V:105:PRO:HD2	1.97	0.44
9:W:10:ILE:HG21	9:W:141:ALA:HB3	1.99	0.44
11:Y:37:ILE:HG23	11:Y:60:GLY:HA2	1.98	0.44
10:J:174:MET:HA	10:X:174:MET:HA	1.98	0.44
9:I:20:VAL:HG13	9:I:118:PRO:HB3	2.00	0.44
13:M:165:ILE:HB	13:M:166:PRO:HD3	2.00	0.44
3:C:201:VAL:HG13	3:C:202:GLN:N	2.33	0.43
9:W:20:VAL:HG13	9:W:118:PRO:HB3	2.00	0.43
3:C:160:GLN:HA	3:C:160:GLN:HE21	1.82	0.43
4:D:160:ASN:HB3	4:D:179:TRP:CE2	2.53	0.43
7:G:78:ILE:N	7:G:79:PRO:CD	2.82	0.43
12:L:13:LEU:CD1	12:L:150:LEU:HD21	2.48	0.43
3:Q:201:VAL:HG13	3:Q:202:GLN:N	2.33	0.43
12:L:4:PRO:O	13:M:104:ARG:NH1	2.46	0.43
12:L:8:ASN:HA	12:L:30:ILE:O	2.19	0.43
9:I:9:GLY:HA3	9:I:41:LYS:HE2	2.02	0.42
1:O:57:MET:HE1	19:U:407:HOH:O	2.19	0.42
7:U:78:ILE:N	7:U:79:PRO:CD	2.82	0.42
9:W:9:GLY:HA3	9:W:41:LYS:HE2	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:Z:8:ASN:HA	12:Z:30:ILE:O	2.19	0.42
6:T:8:ASN:HB2	19:T:301:HOH:O	2.18	0.42
9:I:36:SER:HB2	10:J:126:VAL:HG11	2.01	0.42
2:P:50:LYS:O	2:P:51:VAL:C	2.58	0.42
2:B:161:ALA:HB3	3:C:52:LEU:HD23	2.02	0.41
7:G:149:ASP:HB2	7:G:150:PRO:CD	2.51	0.41
2:B:50:LYS:O	2:B:51:VAL:C	2.59	0.41
7:U:26:THR:HG21	7:U:131:ILE:HD12	2.02	0.41
12:Z:147:MET:N	12:Z:148:PRO:HD2	2.35	0.41
2:B:86:LEU:HB3	2:B:114:LEU:HD21	2.03	0.41
4:D:64:ARG:HG3	19:D:315:HOH:O	2.20	0.41
10:J:93:ARG:NH1	19:J:202:HOH:O	2.53	0.41
11:Y:20:ALA:HB2	11:Y:31:VAL:HG21	2.03	0.41
17:V:301:3BV:H17	9:W:125:LEU:HD22	2.03	0.41
10:X:55:GLN:NE2	19:X:202:HOH:O	2.52	0.41
8:H:168:GLY:O	17:H:301:3BV:H57	2.21	0.41
11:K:20:ALA:HB2	11:K:31:VAL:HG21	2.02	0.41
12:Z:125:PHE:CD2	12:Z:131:TYR:HB3	2.56	0.41
12:Z:13:LEU:CD1	12:Z:150:LEU:HD21	2.48	0.41
2:P:151:ASN:HB2	2:P:152:PRO:CD	2.51	0.41
7:U:149:ASP:HB2	7:U:150:PRO:CD	2.51	0.41
5:S:77:ALA:N	5:S:78:PRO:CD	2.84	0.41
2:B:151:ASN:HB2	2:B:152:PRO:CD	2.50	0.41
5:E:77:ALA:N	5:E:78:PRO:CD	2.84	0.40
7:G:26:THR:HG21	7:G:131:ILE:HD12	2.02	0.40
12:L:147:MET:N	12:L:148:PRO:HD2	2.36	0.40
11:Y:100:MET:CE	11:Y:127:PHE:HB2	2.52	0.40
11:K:100:MET:CE	11:K:127:PHE:HB2	2.52	0.40
8:V:168:GLY:O	17:V:301:3BV:H57	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	A	248/250 (99%)	242 (98%)	6 (2%)	0	100	100
1	O	248/250 (99%)	242 (98%)	6 (2%)	0	100	100
2	B	242/258 (94%)	234 (97%)	6 (2%)	2 (1%)	24	41
2	P	242/258 (94%)	234 (97%)	6 (2%)	2 (1%)	24	41
3	C	238/254 (94%)	230 (97%)	6 (2%)	2 (1%)	24	41
3	Q	238/254 (94%)	231 (97%)	5 (2%)	2 (1%)	24	41
4	D	231/260 (89%)	226 (98%)	5 (2%)	0	100	100
4	R	231/260 (89%)	226 (98%)	5 (2%)	0	100	100
5	E	229/234 (98%)	223 (97%)	6 (3%)	0	100	100
5	S	229/234 (98%)	223 (97%)	6 (3%)	0	100	100
6	F	241/288 (84%)	236 (98%)	5 (2%)	0	100	100
6	T	241/288 (84%)	236 (98%)	5 (2%)	0	100	100
7	G	239/252 (95%)	235 (98%)	4 (2%)	0	100	100
7	U	239/252 (95%)	235 (98%)	4 (2%)	0	100	100
8	H	220/232 (95%)	214 (97%)	6 (3%)	0	100	100
8	V	220/232 (95%)	214 (97%)	6 (3%)	0	100	100
9	I	202/205 (98%)	195 (96%)	7 (4%)	0	100	100
9	W	202/205 (98%)	195 (96%)	7 (4%)	0	100	100
10	J	193/198 (98%)	189 (98%)	3 (2%)	1 (0%)	34	55
10	X	193/198 (98%)	189 (98%)	3 (2%)	1 (0%)	34	55
11	K	210/212 (99%)	207 (99%)	3 (1%)	0	100	100
11	Y	210/212 (99%)	207 (99%)	3 (1%)	0	100	100
12	L	220/222 (99%)	217 (99%)	3 (1%)	0	100	100
12	Z	220/222 (99%)	217 (99%)	3 (1%)	0	100	100
13	M	231/246 (94%)	220 (95%)	10 (4%)	1 (0%)	39	61
13	a	231/246 (94%)	221 (96%)	9 (4%)	1 (0%)	39	61
14	N	194/196 (99%)	188 (97%)	6 (3%)	0	100	100
14	b	194/196 (99%)	188 (97%)	6 (3%)	0	100	100
All	All	6276/6614 (95%)	6114 (97%)	150 (2%)	12 (0%)	52	75

All (12) Ramachandran outliers are listed below:



Mol	Chain	Res	Type
2	B	51	VAL
3	C	202	GLN
2	P	51	VAL
3	Q	202	GLN
3	C	205	ALA
3	Q	205	ALA
2	B	221	ASP
10	J	2	ASP
2	P	221	ASP
10	X	2	ASP
13	M	229	GLY
13	a	229	GLY

### 5.3.2 Protein sidechains ⓘ

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%)	206 (99%)	3 (1%)	74	91
1	O	209/209 (100%)	206 (99%)	3 (1%)	74	91
2	B	203/216 (94%)	197 (97%)	6 (3%)	48	76
2	P	203/216 (94%)	197 (97%)	6 (3%)	48	76
3	C	212/226 (94%)	205 (97%)	7 (3%)	45	73
3	Q	212/226 (94%)	205 (97%)	7 (3%)	45	73
4	D	194/215 (90%)	186 (96%)	8 (4%)	37	63
4	R	194/215 (90%)	186 (96%)	8 (4%)	37	63
5	E	190/193 (98%)	183 (96%)	7 (4%)	41	68
5	S	190/193 (98%)	183 (96%)	7 (4%)	41	68
6	F	201/239 (84%)	193 (96%)	8 (4%)	38	64
6	T	201/239 (84%)	192 (96%)	9 (4%)	34	59
7	G	206/210 (98%)	200 (97%)	6 (3%)	50	77
7	U	206/210 (98%)	200 (97%)	6 (3%)	50	77
8	H	181/190 (95%)	177 (98%)	4 (2%)	60	84

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
8	V	181/190 (95%)	178 (98%)	3 (2%)	68	89
9	I	172/173 (99%)	169 (98%)	3 (2%)	68	89
9	W	172/173 (99%)	169 (98%)	3 (2%)	68	89
10	J	173/175 (99%)	168 (97%)	5 (3%)	50	77
10	X	173/175 (99%)	168 (97%)	5 (3%)	50	77
11	K	169/169 (100%)	164 (97%)	5 (3%)	48	76
11	Y	169/169 (100%)	164 (97%)	5 (3%)	48	76
12	L	185/185 (100%)	180 (97%)	5 (3%)	52	79
12	Z	185/185 (100%)	180 (97%)	5 (3%)	52	79
13	M	199/208 (96%)	193 (97%)	6 (3%)	48	76
13	a	199/208 (96%)	193 (97%)	6 (3%)	48	76
14	N	162/162 (100%)	161 (99%)	1 (1%)	90	97
14	b	162/162 (100%)	161 (99%)	1 (1%)	90	97
All	All	5312/5540 (96%)	5164 (97%)	148 (3%)	51	78

All (148) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	LYS
1	A	61	LEU
1	A	157	PHE
2	B	50	LYS
2	B	55	LEU
2	B	58	GLN
2	B	114	LEU
2	B	184	LYS
2	B	191	LEU
3	C	4	ARG
3	C	38	ASN
3	C	77	ASN
3	C	147	GLN
3	C	160	GLN
3	C	169	VAL
3	C	180	LYS
4	D	99	ILE
4	D	125	LEU
4	D	176	LEU

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Mol	Chain	Res	Type
4	D	193	LEU
4	D	214	ILE
4	D	224	ASP
4	D	236	LYS
4	D	242	GLU
5	E	9	THR
5	E	29	LYS
5	E	54	GLU
5	E	71	LEU
5	E	184	ASN
5	E	188	LEU
5	E	202	ASP
6	F	14	ASP
6	F	123	ASN
6	F	172	LEU
6	F	181	GLU
6	F	203	ASN
6	F	206	LYS
6	F	214	TRP
6	F	221	ASN
7	G	83	ASN
7	G	115	LEU
7	G	125	MET
7	G	166	GLN
7	G	181	LYS
7	G	235	ARG
8	H	30	ASN
8	H	56	THR
8	H	68	LEU
8	H	196	ARG
9	I	37	ASN
9	I	171	LEU
9	I	182	TRP
10	J	23	ARG
10	J	75	LEU
10	J	78	GLN
10	J	144	LEU
10	J	174	MET
11	K	4	LEU
11	K	9	GLN
11	K	35	ILE
11	K	73	ARG

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Mol	Chain	Res	Type
11	K	106	ARG
12	L	1	GLN
12	L	23	LEU
12	L	49	ASN
12	L	136	CYS
12	L	167	LYS
13	M	43	ILE
13	M	48	ASN
13	M	70	LEU
13	M	104	ARG
13	M	161	ARG
13	M	187	ARG
14	N	22	THR
1	O	17	LYS
1	O	61	LEU
1	O	157	PHE
2	P	50	LYS
2	P	55	LEU
2	P	58	GLN
2	P	114	LEU
2	P	184	LYS
2	P	191	LEU
3	Q	4	ARG
3	Q	38	ASN
3	Q	77	ASN
3	Q	147	GLN
3	Q	160	GLN
3	Q	169	VAL
3	Q	180	LYS
4	R	99	ILE
4	R	125	LEU
4	R	176	LEU
4	R	193	LEU
4	R	214	ILE
4	R	224	ASP
4	R	236	LYS
4	R	242	GLU
5	S	9	THR
5	S	29	LYS
5	S	54	GLU
5	S	71	LEU
5	S	184	ASN

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Mol	Chain	Res	Type
5	S	188	LEU
5	S	202	ASP
6	T	14	ASP
6	T	117	GLN
6	T	123	ASN
6	T	172	LEU
6	T	181	GLU
6	T	203	ASN
6	T	206	LYS
6	T	214	TRP
6	T	221	ASN
7	U	83	ASN
7	U	115	LEU
7	U	125	MET
7	U	166	GLN
7	U	181	LYS
7	U	235	ARG
8	V	30	ASN
8	V	56	THR
8	V	196	ARG
9	W	37	ASN
9	W	171	LEU
9	W	182	TRP
10	X	23	ARG
10	X	75	LEU
10	X	78	GLN
10	X	144	LEU
10	X	174	MET
11	Y	4	LEU
11	Y	9	GLN
11	Y	35	ILE
11	Y	73	ARG
11	Y	106	ARG
12	Z	1	GLN
12	Z	23	LEU
12	Z	49	ASN
12	Z	136	CYS
12	Z	167	LYS
13	a	43	ILE
13	a	48	ASN
13	a	70	LEU
13	a	104	ARG

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Mol	Chain	Res	Type
13	a	161	ARG
13	a	187	ARG
14	b	22	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (76) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	94	HIS
2	B	95	GLN
2	B	119	GLN
2	B	123	GLN
3	C	38	ASN
3	C	147	GLN
3	C	160	GLN
4	D	146	GLN
5	E	68	HIS
5	E	92	ASN
5	E	99	ASN
5	E	116	GLN
5	E	118	ASN
5	E	120	GLN
5	E	184	ASN
6	F	86	ASN
6	F	123	ASN
6	F	191	GLN
6	F	240	GLN
7	G	117	GLN
7	G	121	GLN
8	H	66	HIS
8	H	116	HIS
10	J	55	GLN
11	K	85	ASN
11	K	176	ASN
12	L	3	ASN
12	L	70	ASN
12	L	76	HIS
12	L	79	HIS
12	L	158	ASN
13	M	48	ASN
13	M	102	GLN
13	M	108	ASN
13	M	179	ASN

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Mol	Chain	Res	Type
13	M	194	ASN
13	M	213	GLN
1	O	94	HIS
2	P	20	GLN
2	P	119	GLN
2	P	123	GLN
3	Q	38	ASN
3	Q	147	GLN
3	Q	160	GLN
4	R	91	HIS
4	R	146	GLN
5	S	68	HIS
5	S	92	ASN
5	S	99	ASN
5	S	116	GLN
5	S	118	ASN
5	S	120	GLN
5	S	184	ASN
6	T	86	ASN
6	T	117	GLN
6	T	123	ASN
6	T	191	GLN
6	T	240	GLN
7	U	83	ASN
7	U	117	GLN
7	U	121	GLN
8	V	116	HIS
9	W	203	GLN
10	X	55	GLN
11	Y	85	ASN
11	Y	176	ASN
12	Z	3	ASN
12	Z	70	ASN
12	Z	79	HIS
12	Z	158	ASN
13	a	48	ASN
13	a	102	GLN
13	a	108	ASN
13	a	179	ASN
13	a	194	ASN
13	a	213	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 ⓘ

Of 21 ligands modelled in this entry, 11 are monoatomic - leaving 10 for Mogul analysis.

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$
17	3BV	H	301	8	54,54,54	1.41	5 (9%)	67,71,71	1.70	9 (13%)
18	MES	H	302	-	12,12,12	2.07	1 (8%)	15,16,16	2.00	2 (13%)
17	3BV	K	301	11	54,54,54	1.82	6 (11%)	67,71,71	1.83	14 (20%)
18	MES	K	303	-	12,12,12	2.08	1 (8%)	15,16,16	1.55	3 (20%)
17	3BV	N	201	14	54,54,54	1.56	4 (7%)	67,71,71	1.54	8 (11%)
17	3BV	V	301	8	54,54,54	1.45	6 (11%)	67,71,71	1.49	7 (10%)
18	MES	V	303	-	12,12,12	2.08	1 (8%)	15,16,16	2.08	3 (20%)
17	3BV	Y	301	11	54,54,54	1.76	7 (12%)	67,71,71	1.77	10 (14%)
18	MES	Y	303	-	12,12,12	2.12	1 (8%)	15,16,16	1.57	3 (20%)
17	3BV	b	201	14	54,54,54	1.42	4 (7%)	67,71,71	1.66	10 (14%)

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
17	3BV	H	301	8	-	0/59/67/67	0/3/3/3
18	MES	H	302	-	-	0/6/14/14	0/1/1/1
17	3BV	K	301	11	-	0/59/67/67	0/3/3/3
18	MES	K	303	-	-	0/6/14/14	0/1/1/1
17	3BV	N	201	14	-	0/59/67/67	0/3/3/3
17	3BV	V	301	8	-	0/59/67/67	0/3/3/3
18	MES	V	303	-	-	0/6/14/14	0/1/1/1
17	3BV	Y	301	11	-	0/59/67/67	0/3/3/3
18	MES	Y	303	-	-	0/6/14/14	0/1/1/1
17	3BV	b	201	14	-	0/59/67/67	0/3/3/3

All (36) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
18	Y	303	MES	C8-S	-6.97	1.67	1.77
18	V	303	MES	C8-S	-6.87	1.67	1.77
18	H	302	MES	C8-S	-6.84	1.67	1.77
18	K	303	MES	C8-S	-6.83	1.67	1.77
17	K	301	3BV	C32-C33	-6.67	1.34	1.51
17	H	301	3BV	C32-C33	-6.32	1.35	1.51
17	N	201	3BV	C32-C33	-6.09	1.36	1.51
17	b	201	3BV	C32-C33	-5.84	1.36	1.51
17	Y	301	3BV	C32-C33	-5.78	1.37	1.51
17	V	301	3BV	C32-C33	-5.75	1.37	1.51
17	Y	301	3BV	C59-C51	-5.63	1.46	1.52
17	K	301	3BV	C59-C51	-5.59	1.46	1.52
17	b	201	3BV	C13-C14	-4.85	1.37	1.51
17	Y	301	3BV	C13-C14	-4.77	1.37	1.51
17	V	301	3BV	C13-C14	-4.40	1.38	1.51
17	K	301	3BV	C13-C14	-4.30	1.39	1.51
17	H	301	3BV	C13-C14	-4.26	1.39	1.51
17	N	201	3BV	C13-C14	-4.26	1.39	1.51
17	H	301	3BV	C59-C51	-3.65	1.48	1.52
17	K	301	3BV	O48-C47	-3.63	1.34	1.43
17	V	301	3BV	O48-C47	-3.49	1.34	1.43
17	V	301	3BV	C59-C51	-3.21	1.49	1.52
17	N	201	3BV	O48-C47	-3.05	1.35	1.43
17	b	201	3BV	O48-C47	-2.64	1.36	1.43
17	K	301	3BV	C23-C28	-2.33	1.46	1.52
17	H	301	3BV	O48-C47	-2.15	1.37	1.43
17	Y	301	3BV	O48-C47	-2.15	1.37	1.43
17	H	301	3BV	C8-N10	-2.14	1.29	1.34
17	Y	301	3BV	C39-N41	-2.10	1.29	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
17	Y	301	3BV	C8-N10	-2.02	1.30	1.34
17	V	301	3BV	C7-N4	2.17	1.51	1.46
17	V	301	3BV	C51-C47	2.75	1.58	1.53
17	b	201	3BV	C51-C47	3.97	1.61	1.53
17	Y	301	3BV	C51-C47	5.37	1.63	1.53
17	K	301	3BV	C51-C47	5.69	1.64	1.53
17	N	201	3BV	C51-C47	6.11	1.65	1.53

All (69) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
17	H	301	3BV	C58-C51-C59	-8.55	98.48	109.73
17	K	301	3BV	C43-C42-N41	-8.33	98.58	109.98
17	Y	301	3BV	C43-C42-N41	-7.94	99.10	109.98
17	V	301	3BV	C58-C51-C59	-6.94	100.61	109.73
17	N	201	3BV	C43-C42-N41	-6.69	100.82	109.98
17	H	301	3BV	C43-C42-N41	-6.39	101.24	109.98
17	b	201	3BV	C43-C42-N41	-6.34	101.30	109.98
17	N	201	3BV	C58-C51-C59	-5.95	101.91	109.73
17	K	301	3BV	C58-C51-C59	-5.72	102.21	109.73
17	V	301	3BV	C43-C42-N41	-5.11	102.98	109.98
17	Y	301	3BV	C58-C51-C59	-4.92	103.26	109.73
17	b	201	3BV	C7-N4-C3	-4.59	104.41	111.14
17	b	201	3BV	C7-N4-C5	-4.52	104.50	111.14
17	V	301	3BV	O48-C47-C42	-3.93	101.88	109.59
17	b	201	3BV	C58-C51-C59	-3.34	105.34	109.73
17	N	201	3BV	C20-C11-N10	-3.08	102.57	111.28
17	K	301	3BV	O48-C47-C42	-2.96	103.78	109.59
17	b	201	3BV	C20-C11-N10	-2.76	103.48	111.28
17	Y	301	3BV	O48-C47-C42	-2.73	104.24	109.59
17	K	301	3BV	C33-C32-C31	-2.65	105.61	113.44
17	Y	301	3BV	O60-C59-C51	-2.61	105.70	111.16
17	H	301	3BV	O60-C59-C51	-2.61	105.71	111.16
17	H	301	3BV	O48-C47-C42	-2.57	104.55	109.59
17	N	201	3BV	C7-N4-C3	-2.48	107.50	111.14
17	K	301	3BV	O60-C59-C51	-2.47	105.99	111.16
17	b	201	3BV	C12-C13-C14	-2.42	104.25	113.18
17	K	301	3BV	C36-C37-C38	-2.39	116.88	120.20
17	K	301	3BV	C24-C23-C28	-2.26	104.78	110.58
17	K	301	3BV	O1-C2-C3	-2.18	106.79	111.83
17	N	201	3BV	O1-C2-C3	-2.15	106.85	111.83
17	H	301	3BV	C12-C11-N10	-2.14	106.76	110.81

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
17	H	301	3BV	C12-C11-C20	-2.12	105.15	110.34
17	N	201	3BV	O1-C6-C5	-2.07	107.04	111.83
17	Y	301	3BV	C25-C24-C23	-2.06	109.29	115.49
17	K	301	3BV	C39-C31-N30	-2.05	105.47	111.28
17	Y	301	3BV	C39-C31-N30	-2.02	105.55	111.28
17	K	301	3BV	C25-C24-C23	-2.01	109.44	115.49
17	V	301	3BV	C12-C11-C20	-2.01	105.42	110.34
17	H	301	3BV	C3-N4-C5	2.06	113.47	108.87
17	b	201	3BV	C6-C5-N4	2.13	113.36	110.11
18	V	303	MES	O3S-S-C8	2.17	109.50	104.99
17	V	301	3BV	C11-N10-C8	2.19	126.47	121.72
17	H	301	3BV	C7-N4-C5	2.21	114.37	111.14
17	Y	301	3BV	C23-N22-C20	2.26	126.72	121.66
17	b	201	3BV	C2-C3-N4	2.32	113.66	110.11
17	K	301	3BV	C44-C43-C42	2.44	120.72	115.64
18	Y	303	MES	O3S-S-C8	2.64	110.47	104.99
18	K	303	MES	O3S-S-C8	2.65	110.50	104.99
17	Y	301	3BV	C12-C11-C20	2.71	116.99	110.34
17	Y	301	3BV	C3-N4-C5	2.73	114.99	108.87
17	H	301	3BV	C7-N4-C3	2.74	115.15	111.14
18	Y	303	MES	O1S-S-C8	2.83	108.87	106.87
17	b	201	3BV	C47-C42-N41	2.89	116.25	110.06
17	K	301	3BV	C3-N4-C5	2.95	115.48	108.87
17	N	201	3BV	C44-C43-C42	3.01	121.91	115.64
17	N	201	3BV	C47-C42-N41	3.08	116.66	110.06
18	K	303	MES	O1S-S-C8	3.10	109.06	106.87
17	K	301	3BV	C12-C11-C20	3.14	118.05	110.34
17	V	301	3BV	C6-C5-N4	3.14	114.92	110.11
18	Y	303	MES	O2S-S-C8	3.18	109.12	106.87
18	K	303	MES	O2S-S-C8	3.28	109.19	106.87
17	K	301	3BV	C7-N4-C5	3.49	116.25	111.14
18	H	302	MES	O1S-S-C8	3.65	109.45	106.87
17	b	201	3BV	C3-N4-C5	3.66	117.08	108.87
17	V	301	3BV	C7-N4-C3	3.77	116.66	111.14
18	V	303	MES	O2S-S-C8	3.96	109.66	106.87
17	Y	301	3BV	C7-N4-C5	5.35	118.98	111.14
18	H	302	MES	O2S-S-C8	5.91	111.05	106.87
18	V	303	MES	O1S-S-C8	5.98	111.09	106.87

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

7 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
17	H	301	3BV	2	0
17	K	301	3BV	4	0
18	K	303	MES	1	0
17	N	201	3BV	1	0
17	V	301	3BV	4	0
18	V	303	MES	1	0
17	Y	301	3BV	2	0

## 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 ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	250/250 (100%)	0.08	13 (5%)	31 35	29, 45, 85, 136	0
1	O	250/250 (100%)	0.18	16 (6%)	23 25	35, 54, 100, 140	0
2	B	244/258 (94%)	0.31	20 (8%)	14 15	30, 52, 99, 152	0
2	P	244/258 (94%)	0.33	20 (8%)	14 15	33, 55, 104, 153	0
3	C	240/254 (94%)	0.39	29 (12%)	6 5	30, 55, 120, 160	0
3	Q	240/254 (94%)	0.82	39 (16%)	2 2	38, 71, 163, 206	0
4	D	235/260 (90%)	0.12	8 (3%)	49 54	36, 56, 87, 140	0
4	R	235/260 (90%)	0.37	16 (6%)	20 23	44, 64, 106, 154	0
5	E	231/234 (98%)	0.22	13 (5%)	28 31	34, 57, 92, 130	0
5	S	231/234 (98%)	0.56	21 (9%)	11 12	40, 69, 121, 153	0
6	F	243/288 (84%)	0.04	15 (6%)	24 27	30, 51, 104, 129	0
6	T	243/288 (84%)	0.26	11 (4%)	37 42	34, 64, 123, 165	0
7	G	241/252 (95%)	0.06	13 (5%)	29 33	28, 48, 90, 141	0
7	U	241/252 (95%)	0.05	10 (4%)	41 46	30, 51, 85, 129	0
8	H	222/232 (95%)	-0.02	4 (1%)	71 75	30, 44, 80, 116	0
8	V	222/232 (95%)	0.07	5 (2%)	64 67	30, 49, 81, 120	0
9	I	204/205 (99%)	-0.34	1 (0%)	91 92	25, 41, 69, 95	0
9	W	204/205 (99%)	-0.24	3 (1%)	76 79	27, 43, 69, 96	0
10	J	195/198 (98%)	-0.18	4 (2%)	67 71	26, 42, 71, 110	0
10	X	195/198 (98%)	-0.09	4 (2%)	67 71	28, 45, 75, 121	0
11	K	212/212 (100%)	-0.21	3 (1%)	78 80	28, 43, 73, 90	0
11	Y	212/212 (100%)	-0.11	2 (0%)	85 88	30, 45, 76, 101	0
12	L	222/222 (100%)	-0.17	2 (0%)	85 88	25, 44, 75, 100	0
12	Z	222/222 (100%)	-0.15	4 (1%)	71 75	27, 45, 77, 100	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
13	M	233/246 (94%)	-0.12	8 (3%) 49 54	24, 43, 71, 87	0
13	a	233/246 (94%)	-0.17	3 (1%) 79 82	24, 44, 69, 84	0
14	N	196/196 (100%)	-0.26	2 (1%) 84 86	25, 39, 71, 92	0
14	b	196/196 (100%)	-0.20	3 (1%) 76 79	28, 41, 72, 97	0
All	All	6336/6614 (95%)	0.07	292 (4%) 36 41	24, 49, 100, 206	0

All (292) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	219	ALA	14.1
2	P	51	VAL	10.2
1	A	1	MET	10.0
2	P	219	ALA	9.6
2	B	51	VAL	8.5
2	P	218	GLY	8.3
1	O	1	MET	8.2
3	Q	238	LYS	8.0
3	Q	50	LEU	7.9
8	H	222	ASP	7.9
8	H	221	CYS	7.8
8	V	222	ASP	7.8
3	Q	49	THR	7.5
2	B	221	ASP	7.4
5	S	202	ASP	7.3
10	X	1	MET	7.2
3	Q	51	LYS	7.1
2	B	220	ASN	7.0
11	Y	212	GLY	6.9
3	Q	48	SER	6.8
4	R	241	ALA	6.8
13	M	233	ILE	6.7
3	Q	206	LYS	6.6
3	Q	204	GLY	6.5
2	P	220	ASN	6.0
3	Q	202	GLN	6.0
6	F	205	GLU	5.8
2	B	218	GLY	5.8
8	V	221	CYS	5.7
4	D	242	GLU	5.6
10	J	1	MET	5.5
6	T	243	ILE	5.5

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Mol	Chain	Res	Type	RSRZ
3	Q	205	ALA	5.5
3	Q	239	GLN	5.4
1	A	2	THR	5.3
1	O	52	SER	5.3
5	E	233	ILE	5.3
6	F	243	ILE	5.3
10	X	194	ASP	5.2
1	O	250	LEU	5.2
3	C	202	GLN	5.1
3	C	225	GLU	5.0
1	O	249	ALA	5.0
4	D	241	ALA	4.9
9	I	1	SER	4.8
13	a	1	THR	4.8
7	G	242	GLN	4.8
3	Q	223	SER	4.7
7	U	242	GLN	4.7
7	U	222	ASP	4.6
3	Q	240	GLU	4.6
5	S	233	ILE	4.5
3	C	238	LYS	4.5
2	P	221	ASP	4.5
4	R	242	GLU	4.5
3	C	240	GLU	4.4
7	U	2	GLY	4.3
2	P	222	GLY	4.2
10	J	194	ASP	4.2
7	G	240	ALA	4.1
3	C	48	SER	4.1
12	L	165	ASN	4.1
6	T	244	ASN	4.0
5	S	173	ARG	3.9
4	R	125	LEU	3.9
6	T	2	THR	3.9
5	E	201	ARG	3.9
11	Y	106	ARG	3.8
1	A	250	LEU	3.8
13	M	1	THR	3.8
2	P	59	ASP	3.8
6	F	244	ASN	3.8
6	F	2	THR	3.7
1	A	248	GLU	3.7

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Mol	Chain	Res	Type	RSRZ
3	Q	225	GLU	3.6
1	O	53	SER	3.6
5	S	225	ASP	3.6
7	G	188	GLU	3.6
3	Q	203	THR	3.6
3	C	180	LYS	3.6
10	X	193	ASP	3.6
5	S	51	ASN	3.6
3	C	236	GLN	3.6
9	W	1	SER	3.6
3	C	50	LEU	3.5
5	S	180	LYS	3.5
6	T	178	HIS	3.5
13	M	232	LYS	3.5
5	E	122	TYR	3.4
7	G	222	ASP	3.4
3	C	206	LYS	3.4
10	J	193	ASP	3.4
3	C	239	GLN	3.4
6	T	215	CYS	3.4
1	A	50	LYS	3.4
3	C	49	THR	3.4
5	S	122	TYR	3.4
6	F	215	CYS	3.4
1	O	201	GLU	3.3
2	P	52	THR	3.3
13	a	47	ASP	3.3
1	A	249	ALA	3.3
5	S	171	LEU	3.3
7	G	179	LYS	3.3
3	C	175	LYS	3.3
3	Q	229	GLN	3.3
5	S	210	LEU	3.2
6	T	14	ASP	3.2
9	W	133	LYS	3.2
1	A	52	SER	3.2
4	R	113	LEU	3.2
5	S	165	GLN	3.2
6	F	166	GLN	3.2
3	Q	236	GLN	3.2
3	Q	180	LYS	3.1
5	E	202	ASP	3.1

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
3	Q	55	THR	3.1
13	M	47	ASP	3.1
14	b	105	LYS	3.1
3	Q	58	THR	3.1
5	S	54	GLU	3.1
3	Q	59	PRO	3.1
3	Q	60	SER	3.1
3	Q	234	ILE	3.1
3	Q	46	ARG	3.1
6	F	181	GLU	3.0
7	G	3	TYR	3.0
12	Z	165	ASN	3.0
2	P	182	ASP	3.0
3	C	235	GLU	3.0
4	R	201	GLU	3.0
3	Q	52	LEU	3.0
4	D	47	THR	3.0
3	Q	181	GLU	3.0
2	B	232	GLN	3.0
5	S	3	ASN	3.0
13	M	231	GLN	3.0
3	Q	141	ASP	3.0
3	C	60	SER	3.0
2	B	203	SER	2.9
6	T	181	GLU	2.9
1	O	203	GLU	2.9
2	P	244	THR	2.9
1	O	169	VAL	2.9
1	A	201	GLU	2.9
2	B	242	GLY	2.8
2	B	182	ASP	2.8
13	M	216	ASN	2.8
7	U	203	ASP	2.8
2	B	222	GLY	2.8
3	Q	191	LYS	2.8
3	C	1	GLY	2.8
7	U	241	GLU	2.8
2	B	225	TYR	2.8
6	F	201	GLU	2.8
3	Q	237	GLU	2.7
1	A	54	PRO	2.7
3	C	228	ASN	2.7

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
3	Q	27	ARG	2.7
4	R	169	GLU	2.7
8	H	198	GLU	2.7
6	T	241	LYS	2.7
12	Z	1	GLN	2.7
2	B	59	ASP	2.7
14	N	195	GLN	2.7
7	U	188	GLU	2.6
2	P	217	LYS	2.6
2	B	61	SER	2.6
5	S	123	GLY	2.6
5	S	50	ARG	2.6
2	P	203	SER	2.6
3	C	216	ASP	2.6
3	C	47	ARG	2.6
2	P	225	TYR	2.6
5	E	54	GLU	2.6
2	B	62	THR	2.6
4	R	230	GLU	2.6
14	b	195	GLN	2.6
14	b	182	GLY	2.6
4	R	2	ARG	2.6
13	a	204	THR	2.6
5	S	58	TYR	2.6
2	B	60	THR	2.5
3	Q	208	ILE	2.5
4	D	1	ASP	2.5
4	D	125	LEU	2.5
2	P	234	ILE	2.5
4	R	54	ASP	2.5
7	G	2	GLY	2.5
12	Z	106	TYR	2.5
11	K	145	LYS	2.5
7	G	241	GLU	2.5
2	B	54	THR	2.5
2	P	60	THR	2.5
4	R	207	ASN	2.5
11	K	212	GLY	2.5
3	C	27	ARG	2.5
4	D	2	ARG	2.5
4	D	141	ALA	2.5
5	E	123	GLY	2.5

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Mol	Chain	Res	Type	RSRZ
5	E	210	LEU	2.4
1	O	29	LYS	2.4
3	C	181	GLU	2.4
6	F	202	ASP	2.4
7	U	3	TYR	2.4
3	Q	227	ILE	2.4
3	Q	1	GLY	2.4
7	U	183	ASP	2.4
7	G	68	ARG	2.4
3	C	167	LYS	2.4
3	Q	235	GLU	2.4
6	T	205	GLU	2.4
6	T	53	LYS	2.4
4	R	157	TYR	2.4
5	E	173	ARG	2.4
7	G	183	ASP	2.4
12	Z	210	ASP	2.4
13	M	69	ASP	2.4
2	P	177	MET	2.3
5	E	218	ASP	2.3
8	H	219	ASN	2.3
5	S	59	GLN	2.3
3	C	46	ARG	2.3
1	A	203	GLU	2.3
7	U	237	VAL	2.3
9	W	191	LYS	2.3
2	B	235	LYS	2.3
3	Q	187	GLU	2.3
6	F	229	GLY	2.3
2	P	50	LYS	2.3
3	C	37	LYS	2.3
3	C	173	LEU	2.3
8	V	219	ASN	2.3
1	O	231	LYS	2.3
5	E	217	LYS	2.3
3	C	187	GLU	2.3
5	E	227	GLU	2.3
1	A	229	THR	2.3
3	C	169	VAL	2.2
4	R	1	ASP	2.2
6	T	230	ASP	2.2
10	X	151	ASP	2.2

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	A	59	GLU	2.2
2	B	50	LYS	2.2
3	Q	175	LYS	2.2
5	S	52	ALA	2.2
6	F	204	LYS	2.2
3	C	188	GLU	2.2
4	R	203	LYS	2.2
6	F	241	LYS	2.2
5	E	207	VAL	2.2
5	S	207	VAL	2.2
13	M	230	THR	2.2
3	Q	188	GLU	2.2
1	O	50	LYS	2.2
6	F	230	ASP	2.2
12	L	210	ASP	2.2
7	U	153	TYR	2.2
4	D	201	GLU	2.2
5	E	3	ASN	2.2
2	P	15	GLU	2.2
3	Q	142	GLU	2.2
8	V	9	ASN	2.2
11	K	147	ASP	2.2
3	C	51	LYS	2.1
1	O	2	THR	2.1
3	C	59	PRO	2.1
7	G	5	ARG	2.1
5	S	194	GLU	2.1
1	O	229	THR	2.1
3	Q	222	LEU	2.1
5	S	204	SER	2.1
4	R	114	ARG	2.1
7	G	17	TYR	2.1
1	O	40	THR	2.1
4	R	217	GLN	2.1
4	R	175	LEU	2.1
1	O	62	SER	2.1
2	P	93	HIS	2.1
7	G	51	PRO	2.1
8	V	173	VAL	2.1
1	A	58	SER	2.0
14	N	9	LYS	2.0
3	Q	56	ARG	2.0

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Mol	Chain	Res	Type	RSRZ
6	F	206	LYS	2.0
1	O	61	LEU	2.0
2	B	55	LEU	2.0
10	J	195	PHE	2.0
2	B	181	ASP	2.0
6	F	207	ASP	2.0
5	S	218	ASP	2.0
2	P	232	GLN	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
15	MG	Z	301	1/1	0.87	0.37	7.37	58,58,58,58	0
18	MES	Y	303	12/12	0.94	0.31	7.33	33,36,53,56	0
18	MES	V	303	12/12	0.83	0.36	6.78	63,66,82,97	0
18	MES	K	303	12/12	0.93	0.32	6.70	34,37,53,56	0
17	3BV	N	201	52/52	0.83	0.25	5.65	28,43,124,130	0
17	3BV	b	201	52/52	0.85	0.22	4.38	30,44,121,126	0
18	MES	H	302	12/12	0.88	0.32	3.08	61,63,70,82	0
17	3BV	V	301	52/52	0.87	0.22	1.70	43,51,94,99	0
15	MG	I	301	1/1	0.95	0.18	1.66	48,48,48,48	0
17	3BV	Y	301	52/52	0.91	0.17	1.54	31,40,92,97	0
17	3BV	K	301	52/52	0.89	0.18	1.44	29,39,91,97	0
17	3BV	H	301	52/52	0.88	0.19	1.31	41,49,95,99	0
15	MG	N	202	1/1	0.97	0.12	-0.21	37,37,37,37	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
15	MG	G	301	1/1	0.88	0.13	-0.28	49,49,49,49	0
15	MG	K	302	1/1	0.99	0.11	-0.76	32,32,32,32	0
16	CL	b	202	1/1	0.97	0.09	-1.27	60,60,60,60	0
16	CL	N	203	1/1	0.97	0.07	-1.89	50,50,50,50	0
15	MG	V	302	1/1	0.89	0.10	-1.90	39,39,39,39	0
15	MG	Y	302	1/1	0.95	0.06	-3.38	48,48,48,48	0
16	CL	G	302	1/1	0.99	0.03	-	35,35,35,35	0
16	CL	U	301	1/1	0.99	0.09	-	37,37,37,37	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.