



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:50 PM GMT

PDB ID : 4Z1L
Title : Yeast 20S proteasome in complex with belactosin C derivative 3
Authors : Huber, E.M.; Groll, M.
Deposited on : 2015-03-27
Resolution : 3.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

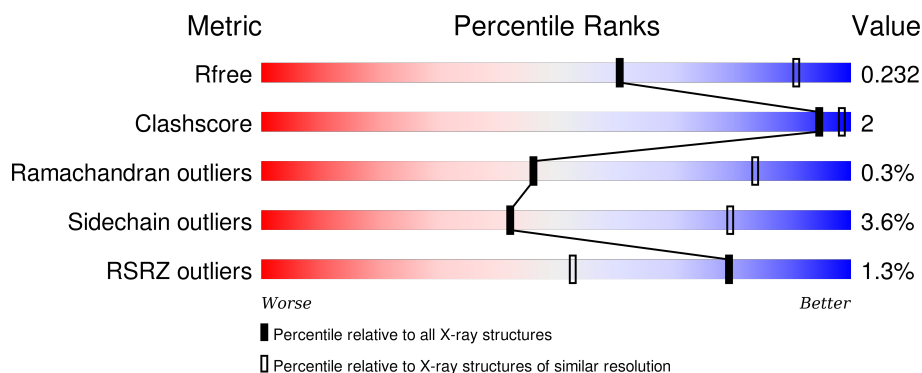
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	1578 (3.00-3.00)
Clashscore	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)
RSRZ outliers	91569	1592 (3.00-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	250	<div> <div>98%</div> <div> <div></div> <div></div> <div></div> <div></div> </div> </div>
1	O	250	<div> <div>2%</div> <div>97%</div> <div> <div></div> <div></div> <div></div> <div></div> </div> </div>
2	B	258	<div> <div>2%</div> <div>86%</div> <div>8%</div> <div>5%</div> <div> <div></div> <div></div> <div></div> <div></div> </div> </div>
2	P	258	<div> <div>2%</div> <div>87%</div> <div>7%</div> <div>5%</div> <div> <div></div> <div></div> <div></div> <div></div> </div> </div>
3	C	254	<div> <div>3%</div> <div>85%</div> <div>8%</div> <div>6%</div> <div> <div></div> <div></div> <div></div> <div></div> </div> </div>

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Mol	Chain	Length	Quality of chain
3	Q	254	
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	I	301	-	-	-	X
15	MG	W	301	-	-	-	X
16	CL	N	203	-	-	-	X
17	4KF	H	301	-	-	-	X
17	4KF	K	301	-	-	-	X
17	4KF	Y	301	-	-	-	X

2 Entry composition

There are 18 unique types of molecules in this entry. The entry contains 49749 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
			1907	1214	320	365	8			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	H	226	Total	C	N	O	S	0	0	0
			1719	1082	298	332	7			
8	V	226	Total	C	N	O	S	0	0	0
			1719	1082	298	332	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	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	229	Total	C	N	O	S	0	0	0
			1790	1133	306	344	7			
13	a	232	Total	C	N	O	S	0	0	0
			1815	1148	311	349	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	0	0
			1	1		
15	K	1	Total	Mg	0	0
			1	1		

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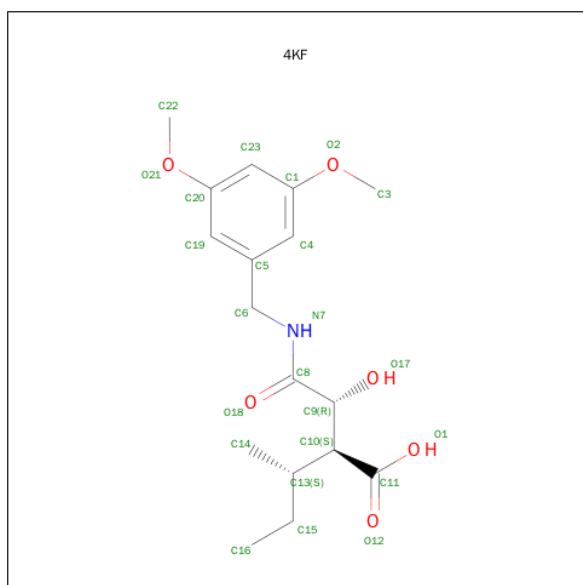
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
15	I	2	Total	Mg	0	0
			2	2		
15	W	1	Total	Mg	0	0
			1	1		
15	Z	1	Total	Mg	0	0
			1	1		
15	N	1	Total	Mg	0	0
			1	1		
15	L	1	Total	Mg	0	0
			1	1		

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

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

- Molecule 17 is (2S,3S)-2-[(1R)-2-[(3,5-dimethoxybenzyl)amino]-1-hydroxy-2-oxoethyl]-3-methylpentanoic acid (three-letter code: 4KF) (formula: C₁₇H₂₅NO₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
17	H	1	Total 23	C 17	N 1	O 5	0	0
17	K	1	Total 23	C 17	N 1	O 5	0	0
17	N	1	Total 23	C 17	N 1	O 5	0	0
17	V	1	Total 23	C 17	N 1	O 5	0	0
17	Y	1	Total 23	C 17	N 1	O 5	0	0
17	b	1	Total 23	C 17	N 1	O 5	0	0

- Molecule 18 is water.

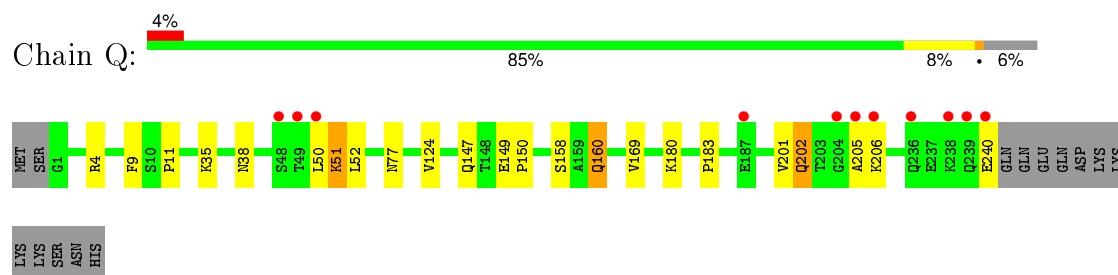
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
18	A	5	Total 5	O 5	0	0
18	B	12	Total 12	O 12	0	0
18	C	8	Total 8	O 8	0	0
18	D	10	Total 10	O 10	0	0
18	E	6	Total 6	O 6	0	0
18	F	11	Total 11	O 11	0	0
18	G	9	Total 9	O 9	0	0
18	H	8	Total 8	O 8	0	0
18	I	7	Total 7	O 7	0	0
18	J	9	Total 9	O 9	0	0
18	K	13	Total 13	O 13	0	0
18	L	21	Total 21	O 21	0	0
18	M	15	Total 15	O 15	0	0
18	N	8	Total 8	O 8	0	0

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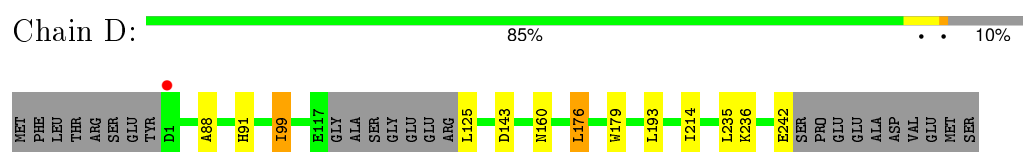
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
18	P	15	Total 15	O 15	0	0
18	Q	16	Total 16	O 16	0	0
18	R	3	Total 3	O 3	0	0
18	S	4	Total 4	O 4	0	0
18	T	8	Total 8	O 8	0	0
18	U	13	Total 13	O 13	0	0
18	V	7	Total 7	O 7	0	0
18	W	9	Total 9	O 9	0	0
18	X	10	Total 10	O 10	0	0
18	Y	14	Total 14	O 14	0	0
18	Z	14	Total 14	O 14	0	0
18	a	8	Total 8	O 8	0	0
18	b	13	Total 13	O 13	0	0

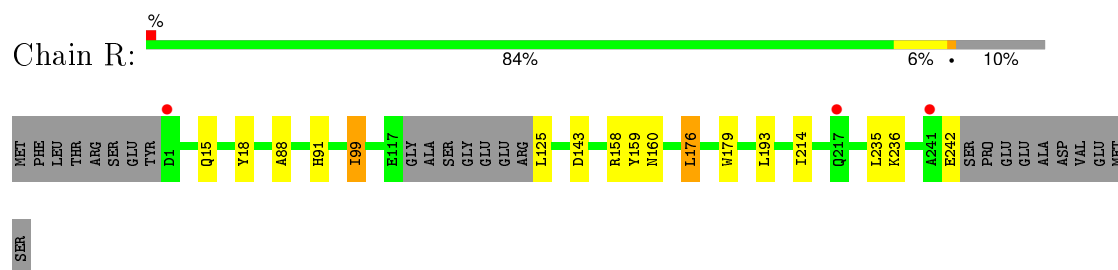
- Molecule 3: Proteasome subunit alpha type-4



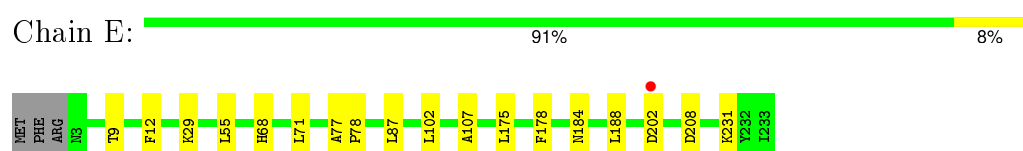
- Molecule 4: Proteasome subunit alpha type-5



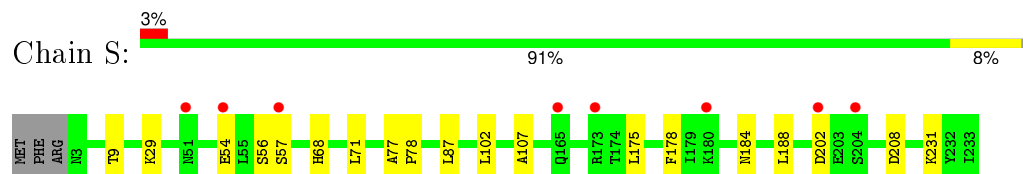
- Molecule 4: Proteasome subunit alpha type-5



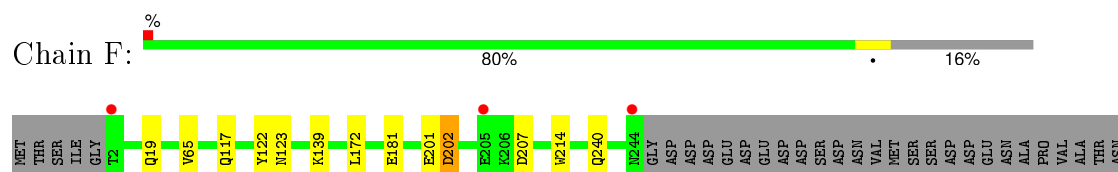
- Molecule 5: Proteasome subunit alpha type-6



- Molecule 5: Proteasome subunit alpha type-6




- Molecule 6: Probable proteasome subunit alpha type-7



ALA
ASN
ALA
THR
THR
GLY
ASP
GLN
GLU
GLY
ASP
ILE
HIS
LEU
GLU


- Molecule 6: Probable proteasome subunit alpha type-7

Chain T:  80% 5% 16%

MET THR SER ILE THR ASP GLY GLN GLU ASP ILE HIS LEU GLU
T2 Q117 Y122 N123 K139 E148 L172 E181 L198 E201 D202 D207 W214 Q240 L243 E244 GLY ASP ASP ASP GLU ASP GLU ASP ASP ASP SER ASP ASN VAL MET SER SER ASP ASP GLU ASN ALA PRO VAL ALA THR ASN ALA


ASN
ALA
THR
THR
ASP
GLY
GLY
ASP
ILE
HIS
LEU
GLU

- Molecule 7: Proteasome subunit alpha type-1

Chain G:  88% 7% 5%


MET SER GLY ALA ALA ALA SER ALA ALA G2 F23 T26 I63 I68 V73 V74 N75 I78 P79 N83 L115 S116 Q117 R122 M125 T133 E208 E215 E235 L236 Q242 ASP

- Molecule 7: Proteasome subunit alpha type-1

Chain U:  88% 8% 4%

MET SER GLY ALA ALA ALA SER ALA ALA G2 P12 F23 T26 I63 V73 V74 N75 I78 P79 N83 L115 S116 Q117 R122 M125 T133 K165 L205 E208 E215 E235 D222 R235 L236 E241 Q242 ASP

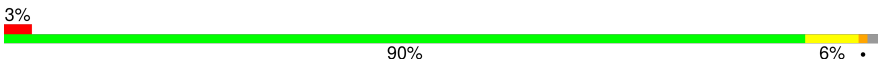
- Molecule 8: Proteasome subunit beta type-2

Chain H:  87% 9% 4%

T1 I14 Q22 N30 H35 A49 T52 T56 I63 L68 P74 L80 G82 I89 V100 A101 D104 P105 F111 S112 I113 H116 T119 S126 L127 R196 E197 E198 T210 C221 D222 I223 Q224 E225 E226 GLN VAL ASP ILE

THR
ALA

- Molecule 8: Proteasome subunit beta type-2

Chain V:  90% 6% 4%

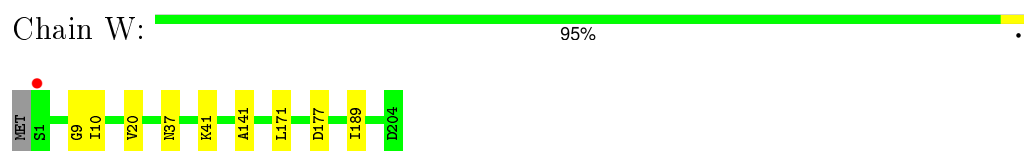
T1 T2 I3 Q22 N30 H35 T52 T56 I63 L68 P74 L80 K84 D104 P105 I113 T119 L127 R196 C221 D222 I223 Q224 E225 E226 GLN VAL ASP ILE THR ALA

- Molecule 9: Proteasome subunit beta type-3

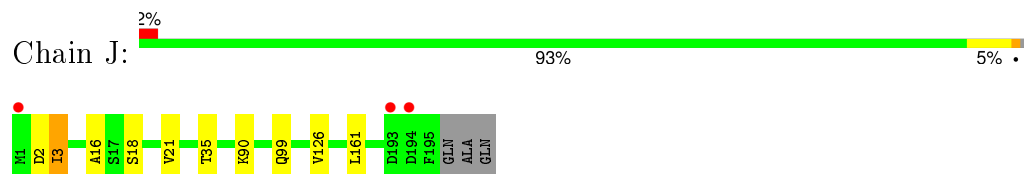
Chain I:  94% 6%

MET S1 G9 I10 V20 S36 N37 K41 P118 A141 S167 L171 D177 I189 D204

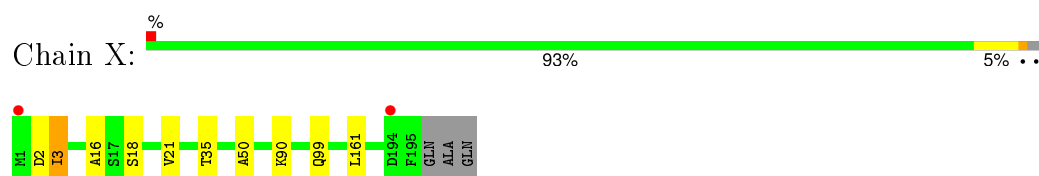
- Molecule 9: Proteasome subunit beta type-3



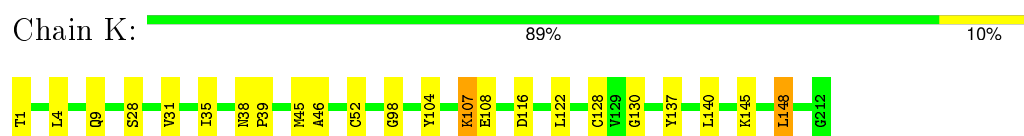
- Molecule 10: Proteasome subunit beta type-4



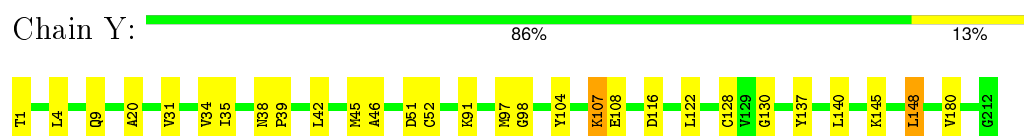
- Molecule 10: Proteasome subunit beta type-4



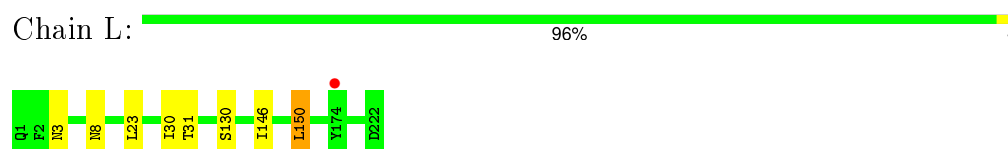
- Molecule 11: Proteasome subunit beta type-5



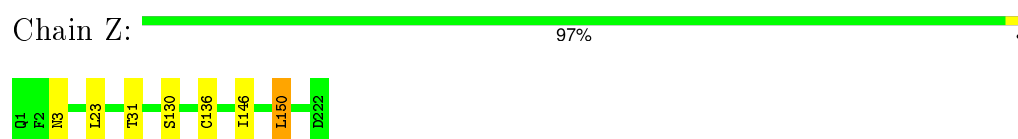
- Molecule 11: Proteasome subunit beta type-5



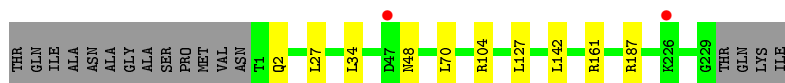
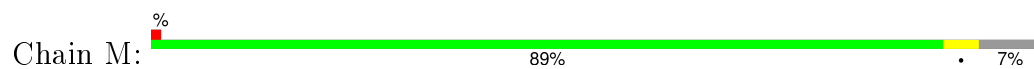
- Molecule 12: Proteasome subunit beta type-6



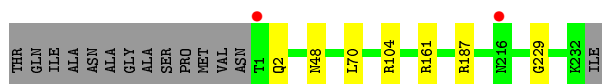
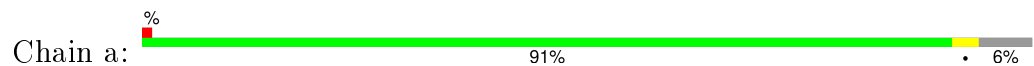
- Molecule 12: Proteasome subunit beta type-6



- 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, α , β , γ	136.09Å 301.30Å 144.58Å 90.00° 113.33° 90.00°	Depositor
Resolution (Å)	15.00 – 3.00 15.00 – 3.00	Depositor EDS
% Data completeness (in resolution range)	97.7 (15.00-3.00) 98.1 (15.00-3.00)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.38 (at 3.01Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.204 , 0.229 0.208 , 0.232	Depositor DCC
R_{free} test set	10364 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	54.8	Xtriage
Anisotropy	0.186	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 42.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	2 of 206885 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	49749	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.37% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 4KF, MG, 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	0.47	0/2642
1	O	0.27	0/1952	0.47	0/2642
2	B	0.28	0/1934	0.50	0/2618
2	P	0.28	0/1934	0.50	0/2618
3	C	0.27	0/1910	0.51	0/2586
3	Q	0.27	0/1910	0.51	0/2586
4	D	0.27	0/1837	0.48	0/2475
4	R	0.26	0/1837	0.48	0/2475
5	E	0.27	0/1800	0.48	0/2433
5	S	0.27	0/1800	0.48	0/2433
6	F	0.27	0/1932	0.46	0/2609
6	T	0.26	0/1932	0.45	0/2609
7	G	0.27	0/1945	0.48	0/2634
7	U	0.27	0/1945	0.48	0/2634
8	H	0.33	0/1750	0.51	0/2373
8	V	0.35	0/1750	0.52	0/2373
9	I	0.28	0/1611	0.49	0/2174
9	W	0.28	0/1611	0.49	0/2174
10	J	0.27	0/1589	0.48	0/2142
10	X	0.27	0/1589	0.48	0/2142
11	K	0.42	0/1681	0.53	0/2274
11	Y	0.33	0/1681	0.53	0/2274
12	L	0.27	0/1795	0.48	0/2420
12	Z	0.28	0/1795	0.48	0/2420
13	M	0.28	0/1821	0.51	0/2470
13	a	0.28	0/1846	0.51	0/2503
14	N	0.29	0/1541	0.48	0/2087
14	b	0.27	0/1541	0.48	0/2087
All	All	0.29	0/50221	0.49	0/67907

There are no bond length outliers.

There are no bond angle outliers.

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	2	0
1	O	1915	0	1929	4	0
2	B	1904	0	1904	12	0
2	P	1904	0	1904	10	0
3	C	1881	0	1895	8	0
3	Q	1881	0	1895	9	0
4	D	1813	0	1797	6	0
4	R	1813	0	1797	10	0
5	E	1773	0	1775	6	0
5	S	1773	0	1775	7	0
6	F	1892	0	1883	4	0
6	T	1892	0	1883	4	0
7	G	1907	0	1901	5	0
7	U	1907	0	1901	6	0
8	H	1719	0	1718	14	0
8	V	1719	0	1718	7	0
9	I	1581	0	1574	7	0
9	W	1581	0	1574	4	0
10	J	1561	0	1569	4	0
10	X	1561	0	1569	4	0
11	K	1644	0	1594	19	0
11	Y	1644	0	1594	19	0
12	L	1757	0	1711	2	0
12	Z	1757	0	1711	1	0
13	M	1790	0	1793	2	0
13	a	1815	0	1821	0	0
14	N	1512	0	1480	4	0
14	b	1512	0	1480	0	0
15	G	1	0	0	0	0
15	I	2	0	0	0	0
15	K	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
15	L	1	0	0	0	0
15	N	1	0	0	0	0
15	W	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	23	0	24	2	0
17	K	23	0	24	7	0
17	N	23	0	24	0	0
17	V	23	0	24	0	0
17	Y	23	0	24	5	0
17	b	23	0	24	0	0
18	A	5	0	0	0	0
18	B	12	0	0	1	0
18	C	8	0	0	0	0
18	D	10	0	0	0	0
18	E	6	0	0	0	0
18	F	11	0	0	1	0
18	G	9	0	0	0	0
18	H	8	0	0	0	0
18	I	7	0	0	0	0
18	J	9	0	0	0	0
18	K	13	0	0	0	0
18	L	21	0	0	0	0
18	M	15	0	0	0	0
18	N	8	0	0	0	0
18	P	15	0	0	1	0
18	Q	16	0	0	1	0
18	R	3	0	0	0	0
18	S	4	0	0	0	0
18	T	8	0	0	1	0
18	U	13	0	0	0	0
18	V	7	0	0	0	0
18	W	9	0	0	0	0
18	X	10	0	0	0	0
18	Y	14	0	0	0	0
18	Z	14	0	0	0	0
18	a	8	0	0	0	0
18	b	13	0	0	0	0
All	All	49749	0	49218	161	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 2.

All (161) 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)
8:H:52:THR:O	8:H:56:THR:OG1	2.03	0.77
8:V:52:THR:O	8:V:56:THR:OG1	2.03	0.77
17:Y:301:4KF:O12	17:Y:301:4KF:H17	1.88	0.72
11:K:45:MET:CE	17:K:301:4KF:H16	2.20	0.71
11:K:45:MET:HE3	17:K:301:4KF:H16	1.76	0.65
11:K:45:MET:HE3	17:K:301:4KF:C16	2.26	0.65
11:Y:34:VAL:CG1	11:Y:42:LEU:HD22	2.26	0.64
2:P:93:HIS:HB3	18:P:301:HOH:O	1.98	0.63
11:Y:51:ASP:HB3	11:Y:97:MET:CE	2.32	0.59
11:K:28:SER:OG	11:K:31:VAL:HG23	2.03	0.59
4:R:159:TYR:CE2	5:S:56:SER:HB3	2.38	0.58
11:Y:51:ASP:HB3	11:Y:97:MET:HE2	1.85	0.58
2:B:93:HIS:HB3	18:B:301:HOH:O	2.03	0.57
17:Y:301:4KF:C11	17:Y:301:4KF:H17	2.33	0.57
8:H:14:ILE:HD12	8:H:101:ALA:CB	2.35	0.56
8:H:14:ILE:HD11	8:H:99:ILE:HG22	1.87	0.55
2:B:12:PHE:H	3:C:17:GLN:HE22	1.55	0.54
8:H:14:ILE:HD12	8:H:101:ALA:HB3	1.89	0.54
11:K:128:CYS:HB2	11:K:137:TYR:CZ	2.43	0.54
4:R:176:LEU:HD11	5:S:54:GLU:HB2	1.91	0.53
10:J:16:ALA:HB2	10:J:161:LEU:HD21	1.91	0.53
11:K:35:ILE:HD11	11:K:45:MET:CG	2.39	0.53
11:Y:128:CYS:HB2	11:Y:137:TYR:CZ	2.45	0.52
6:T:148:GLU:HG2	18:T:302:HOH:O	2.08	0.52
10:X:16:ALA:HB2	10:X:161:LEU:HD21	1.91	0.52
11:Y:1:THR:O	11:Y:130:GLY:HA3	2.09	0.52
7:U:23:PHE:O	7:U:26:THR:HB	2.10	0.52
7:G:23:PHE:O	7:G:26:THR:HB	2.10	0.52
11:K:45:MET:HB3	17:K:301:4KF:C16	2.40	0.52
8:H:113:ILE:HG23	8:H:119:THR:HG22	1.91	0.51
3:C:160:GLN:HA	3:C:160:GLN:HE21	1.75	0.51
3:C:201:VAL:O	3:C:202:GLN:CB	2.57	0.51
7:G:68:ARG:HH12	14:N:36:ARG:HH22	1.58	0.51
3:Q:201:VAL:O	3:Q:202:GLN:CB	2.58	0.51
8:H:92:GLY:HA3	8:H:116:HIS:CD2	2.47	0.50
3:Q:160:GLN:HE21	3:Q:160:GLN:HA	1.75	0.50
14:N:83:LYS:HG3	14:N:119:VAL:CG2	2.42	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:W:9:GLY:HA3	9:W:41:LYS:HE2	1.94	0.50
3:Q:51:LYS:O	3:Q:52:LEU:HB2	2.11	0.50
9:I:9:GLY:HA3	9:I:41:LYS:HE2	1.94	0.50
3:C:51:LYS:O	3:C:52:LEU:HB2	2.10	0.49
8:V:80:LEU:HD13	8:V:119:THR:HG21	1.94	0.49
11:K:45:MET:HE2	17:K:301:4KF:H16	1.95	0.49
8:H:49:ALA:HB2	17:H:301:4KF:H17	1.93	0.49
11:Y:42:LEU:HD12	11:Y:180:VAL:HG22	1.94	0.49
2:P:47:ALA:HB1	2:P:64:LYS:HD2	1.95	0.49
5:E:12:PHE:H	6:F:19:GLN:HE22	1.60	0.48
17:Y:301:4KF:C11	17:Y:301:4KF:C16	2.90	0.48
9:I:36:SER:HB2	10:J:126:VAL:HG11	1.94	0.48
4:R:158:ARG:HB3	5:S:57:SER:HB3	1.95	0.48
1:O:23:TYR:CD1	7:U:12:PRO:HA	2.49	0.48
2:B:47:ALA:HB1	2:B:64:LYS:HD2	1.96	0.48
11:Y:20:ALA:HB2	11:Y:31:VAL:HG21	1.96	0.48
5:S:87:LEU:HD21	5:S:107:ALA:HB1	1.95	0.47
3:C:35:LYS:HG2	3:C:158:SER:O	2.15	0.47
8:H:35:HIS:HB2	8:H:56:THR:HG21	1.97	0.47
11:K:128:CYS:HB2	11:K:137:TYR:CE2	2.49	0.47
11:Y:35:ILE:HD11	11:Y:45:MET:HG2	1.97	0.47
6:F:65:VAL:HG12	18:F:301:HOH:O	2.15	0.47
8:V:35:HIS:HB2	8:V:56:THR:HG21	1.97	0.46
11:Y:1:THR:HG21	17:Y:301:4KF:O12	2.14	0.46
14:N:1:THR:O	14:N:128:GLY:HA3	2.14	0.46
11:K:45:MET:HB3	17:K:301:4KF:H15	1.98	0.46
11:K:45:MET:HG2	11:K:52:CYS:HB3	1.97	0.46
5:E:87:LEU:HD21	5:E:107:ALA:HB1	1.96	0.46
8:H:126:SER:O	8:H:127:LEU:HD13	2.16	0.46
8:V:3:ILE:HD11	8:V:127:LEU:HB3	1.98	0.46
11:Y:1:THR:OG1	17:Y:301:4KF:O17	2.28	0.46
11:Y:52:CYS:SG	11:Y:97:MET:HG3	2.56	0.46
3:Q:52:LEU:N	18:Q:301:HOH:O	2.47	0.46
11:Y:128:CYS:HB2	11:Y:137:TYR:CE2	2.51	0.45
2:P:217:LYS:C	2:P:219:ALA:H	2.19	0.45
10:X:50:ALA:O	11:Y:91:LYS:NZ	2.48	0.45
11:Y:35:ILE:HD11	11:Y:45:MET:CG	2.47	0.45
3:Q:35:LYS:HG2	3:Q:158:SER:O	2.16	0.45
4:D:176:LEU:HD22	5:E:55:LEU:CD2	2.47	0.45
2:B:217:LYS:C	2:B:219:ALA:H	2.19	0.45
4:D:91:HIS:HB3	4:D:99:ILE:CG2	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:W:20:VAL:HG23	9:W:189:ILE:HB	1.99	0.45
3:Q:9:PHE:H	4:R:15:GLN:HE22	1.65	0.44
1:O:122:THR:HG22	2:P:128:ARG:HH21	1.83	0.44
4:D:88:ALA:HA	4:D:99:ILE:HG21	2.00	0.44
10:X:21:VAL:HG11	11:Y:122:LEU:HD11	2.00	0.44
8:V:63:ILE:HG23	8:V:74:PRO:HB3	2.00	0.44
9:I:141:ALA:HB2	9:I:177:ASP:HB2	2.00	0.44
4:R:91:HIS:HB3	4:R:99:ILE:CG2	2.48	0.44
8:H:104:ASP:HB2	8:H:105:PRO:HD2	1.99	0.44
8:V:104:ASP:HB2	8:V:105:PRO:HD2	2.00	0.44
5:S:68:HIS:HE1	5:S:102:LEU:O	2.00	0.44
4:R:160:ASN:HB3	4:R:179:TRP:CE2	2.53	0.43
5:E:68:HIS:HE1	5:E:102:LEU:O	2.01	0.43
4:R:88:ALA:HA	4:R:99:ILE:HG21	2.00	0.43
6:F:202:ASP:OD1	6:F:202:ASP:N	2.51	0.43
11:K:1:THR:O	11:K:130:GLY:HA3	2.18	0.43
9:I:10:ILE:HG21	9:I:141:ALA:HB3	2.00	0.43
8:V:84:LYS:HA	8:V:113:ILE:HD11	2.00	0.43
11:K:107:LYS:HG3	11:K:108:GLU:HG3	2.00	0.43
2:P:151:ASN:HB2	2:P:152:PRO:HD2	2.01	0.43
6:T:202:ASP:N	6:T:202:ASP:OD1	2.52	0.43
7:U:73:VAL:HG12	7:U:133:THR:HB	2.00	0.43
9:W:141:ALA:HB2	9:W:177:ASP:HB2	2.00	0.43
11:K:45:MET:HE3	17:K:301:4KF:H15	1.99	0.43
2:B:124:HIS:HB3	3:C:124:VAL:HG12	2.01	0.43
11:Y:107:LYS:HG3	11:Y:108:GLU:HG3	2.01	0.43
4:D:91:HIS:HB3	4:D:99:ILE:HG22	2.01	0.43
4:D:160:ASN:HB3	4:D:179:TRP:CE2	2.54	0.43
8:H:63:ILE:HG23	8:H:74:PRO:HB3	2.01	0.43
10:J:21:VAL:HG11	11:K:122:LEU:HD11	2.00	0.43
7:G:73:VAL:HG12	7:G:133:THR:HB	2.00	0.43
9:I:20:VAL:HG23	9:I:189:ILE:HB	1.99	0.43
2:B:50:LYS:HD3	2:B:50:LYS:HA	1.87	0.42
3:Q:149:GLU:HB2	3:Q:150:PRO:HD2	2.01	0.42
5:E:175:LEU:HA	5:E:178:PHE:CE2	2.54	0.42
12:Z:146:ILE:HG22	12:Z:150:LEU:HD22	2.01	0.42
9:W:10:ILE:HG21	9:W:141:ALA:HB3	2.01	0.42
11:K:35:ILE:CD1	11:K:45:MET:CG	2.98	0.42
1:A:1:MET:HG3	6:F:122:TYR:CZ	2.54	0.42
5:S:175:LEU:HA	5:S:178:PHE:CE2	2.54	0.42
4:R:91:HIS:HB3	4:R:99:ILE:HG22	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:14:PRO:HA	3:C:20:TYR:CD1	2.54	0.42
12:L:146:ILE:HG22	12:L:150:LEU:HD22	2.01	0.42
2:B:151:ASN:HB2	2:B:152:PRO:CD	2.50	0.42
3:C:149:GLU:HB2	3:C:150:PRO:HD2	2.02	0.42
14:N:36:ARG:HG3	14:N:42:TRP:CE2	2.54	0.42
2:B:151:ASN:HB2	2:B:152:PRO:HD2	2.00	0.42
10:J:3:ILE:HG23	10:J:18:SER:HB3	2.01	0.42
8:H:80:LEU:HD13	8:H:111:PHE:CG	2.55	0.42
8:H:210:THR:HG21	9:I:167:SER:HB3	2.00	0.42
7:U:78:ILE:N	7:U:79:PRO:CD	2.83	0.42
13:M:127:LEU:HG	13:M:142:LEU:HD12	2.02	0.42
10:X:3:ILE:HG23	10:X:18:SER:HB3	2.01	0.42
1:A:122:THR:HG22	2:B:128:ARG:HH21	1.85	0.41
11:K:46:ALA:HB3	11:K:98:GLY:O	2.19	0.41
11:Y:38:ASN:HB2	11:Y:39:PRO:CD	2.50	0.41
2:P:151:ASN:HB2	2:P:152:PRO:CD	2.50	0.41
2:P:219:ALA:HB2	2:P:225:TYR:HB2	2.01	0.41
2:B:219:ALA:HB2	2:B:225:TYR:HB2	2.02	0.41
2:B:50:LYS:O	2:B:51:VAL:C	2.58	0.41
11:Y:46:ALA:HB3	11:Y:98:GLY:O	2.19	0.41
7:G:78:ILE:N	7:G:79:PRO:CD	2.83	0.41
11:K:38:ASN:HB2	11:K:39:PRO:CD	2.50	0.41
7:G:63:ILE:HD12	7:G:215:GLU:HG2	2.03	0.41
2:P:50:LYS:O	2:P:51:VAL:C	2.58	0.41
11:K:145:LYS:HB2	11:K:148:LEU:HD13	2.03	0.41
1:O:1:MET:HG3	6:T:122:TYR:CZ	2.56	0.41
5:S:77:ALA:N	5:S:78:PRO:CD	2.84	0.41
3:Q:11:PRO:HA	4:R:18:TYR:CD1	2.56	0.41
17:H:301:4KF:H4	17:H:301:4KF:H16	1.87	0.40
11:Y:145:LYS:HB2	11:Y:148:LEU:HD13	2.03	0.40
5:E:77:ALA:N	5:E:78:PRO:CD	2.84	0.40
4:D:91:HIS:CD2	4:D:99:ILE:HG22	2.57	0.40
7:U:63:ILE:HD12	7:U:215:GLU:HG2	2.03	0.40
9:I:20:VAL:HG13	9:I:118:PRO:HB3	2.03	0.40
12:L:8:ASN:HA	12:L:30:ILE:O	2.22	0.40
6:T:198:LEU:HD12	6:T:243:ILE:HG22	2.03	0.40
2:P:124:HIS:HB3	3:Q:124:VAL:HG12	2.04	0.40
7:U:165:LYS:HD2	7:U:205:LEU:HD22	2.03	0.40
4:R:91:HIS:CD2	4:R:99:ILE:HG22	2.57	0.40
8:H:104:ASP:HB2	8:H:105:PRO:CD	2.51	0.40
1:O:161:ALA:O	2:P:55:LEU:HD23	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:M:27:LEU:HD21	13:M:34:LEU:HD22	2.02	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%)	238 (96%)	9 (4%)	1 (0%)	39	80
1	O	248/250 (99%)	238 (96%)	9 (4%)	1 (0%)	39	80
2	B	242/258 (94%)	235 (97%)	3 (1%)	4 (2%)	11	46
2	P	242/258 (94%)	235 (97%)	3 (1%)	4 (2%)	11	46
3	C	238/254 (94%)	233 (98%)	2 (1%)	3 (1%)	15	53
3	Q	238/254 (94%)	233 (98%)	2 (1%)	3 (1%)	15	53
4	D	231/260 (89%)	228 (99%)	3 (1%)	0	100	100
4	R	231/260 (89%)	228 (99%)	3 (1%)	0	100	100
5	E	229/234 (98%)	222 (97%)	7 (3%)	0	100	100
5	S	229/234 (98%)	222 (97%)	7 (3%)	0	100	100
6	F	241/288 (84%)	237 (98%)	4 (2%)	0	100	100
6	T	241/288 (84%)	236 (98%)	5 (2%)	0	100	100
7	G	239/252 (95%)	237 (99%)	2 (1%)	0	100	100
7	U	239/252 (95%)	237 (99%)	2 (1%)	0	100	100
8	H	224/232 (97%)	219 (98%)	5 (2%)	0	100	100
8	V	224/232 (97%)	219 (98%)	5 (2%)	0	100	100
9	I	202/205 (98%)	194 (96%)	8 (4%)	0	100	100
9	W	202/205 (98%)	194 (96%)	8 (4%)	0	100	100
10	J	193/198 (98%)	188 (97%)	5 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
10	X	193/198 (98%)	188 (97%)	5 (3%)	0	100	100
11	K	210/212 (99%)	205 (98%)	5 (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%)	214 (97%)	6 (3%)	0	100	100
13	M	227/246 (92%)	221 (97%)	6 (3%)	0	100	100
13	a	230/246 (94%)	223 (97%)	6 (3%)	1 (0%)	39	80
14	N	194/196 (99%)	190 (98%)	4 (2%)	0	100	100
14	b	194/196 (99%)	190 (98%)	4 (2%)	0	100	100
All	All	6279/6614 (95%)	6125 (98%)	137 (2%)	17 (0%)	46	84

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	51	VAL
2	B	221	ASP
3	C	202	GLN
2	P	51	VAL
2	P	221	ASP
3	Q	202	GLN
1	A	2	THR
2	B	218	GLY
2	B	220	ASN
1	O	2	THR
2	P	218	GLY
2	P	220	ASN
3	C	205	ALA
3	Q	205	ALA
3	C	183	PRO
3	Q	183	PRO
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%)	205 (98%)	4 (2%)	65	90
1	O	209/209 (100%)	205 (98%)	4 (2%)	65	90
2	B	203/216 (94%)	197 (97%)	6 (3%)	48	83
2	P	203/216 (94%)	197 (97%)	6 (3%)	48	83
3	C	212/226 (94%)	201 (95%)	11 (5%)	29	68
3	Q	212/226 (94%)	201 (95%)	11 (5%)	29	68
4	D	194/215 (90%)	185 (95%)	9 (5%)	33	73
4	R	194/215 (90%)	185 (95%)	9 (5%)	33	73
5	E	190/193 (98%)	182 (96%)	8 (4%)	36	76
5	S	190/193 (98%)	182 (96%)	8 (4%)	36	76
6	F	201/239 (84%)	191 (95%)	10 (5%)	30	70
6	T	201/239 (84%)	191 (95%)	10 (5%)	30	70
7	G	206/210 (98%)	196 (95%)	10 (5%)	31	71
7	U	206/210 (98%)	196 (95%)	10 (5%)	31	71
8	H	185/190 (97%)	178 (96%)	7 (4%)	40	78
8	V	185/190 (97%)	178 (96%)	7 (4%)	40	78
9	I	172/173 (99%)	170 (99%)	2 (1%)	78	94
9	W	172/173 (99%)	170 (99%)	2 (1%)	78	94
10	J	173/175 (99%)	168 (97%)	5 (3%)	50	84
10	X	173/175 (99%)	168 (97%)	5 (3%)	50	84
11	K	169/169 (100%)	162 (96%)	7 (4%)	37	76
11	Y	169/169 (100%)	162 (96%)	7 (4%)	37	76
12	L	185/185 (100%)	180 (97%)	5 (3%)	52	85
12	Z	185/185 (100%)	179 (97%)	6 (3%)	46	82
13	M	195/208 (94%)	189 (97%)	6 (3%)	47	83
13	a	198/208 (95%)	192 (97%)	6 (3%)	48	83
14	N	162/162 (100%)	158 (98%)	4 (2%)	55	86
14	b	162/162 (100%)	158 (98%)	4 (2%)	55	86
All	All	5315/5540 (96%)	5126 (96%)	189 (4%)	42	79

All (189) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	62	SER
1	A	122	THR
1	A	157	PHE
1	A	250	LEU
2	B	50	LYS
2	B	55	LEU
2	B	79	LEU
2	B	113	ARG
2	B	191	LEU
2	B	238	LEU
3	C	4	ARG
3	C	38	ASN
3	C	50	LEU
3	C	51	LYS
3	C	77	ASN
3	C	147	GLN
3	C	160	GLN
3	C	169	VAL
3	C	180	LYS
3	C	206	LYS
3	C	240	GLU
4	D	99	ILE
4	D	125	LEU
4	D	143	ASP
4	D	176	LEU
4	D	193	LEU
4	D	214	ILE
4	D	235	LEU
4	D	236	LYS
4	D	242	GLU
5	E	9	THR
5	E	29	LYS
5	E	71	LEU
5	E	184	ASN
5	E	188	LEU
5	E	202	ASP
5	E	208	ASP
5	E	231	LYS
6	F	117	GLN
6	F	123	ASN
6	F	139	LYS
6	F	172	LEU
6	F	181	GLU

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Mol	Chain	Res	Type
6	F	201	GLU
6	F	202	ASP
6	F	207	ASP
6	F	214	TRP
6	F	240	GLN
7	G	26	THR
7	G	75	ASN
7	G	83	ASN
7	G	115	LEU
7	G	117	GLN
7	G	122	ARG
7	G	125	MET
7	G	208	GLU
7	G	235	ARG
7	G	236	LEU
8	H	14	ILE
8	H	22	GLN
8	H	30	ASN
8	H	56	THR
8	H	68	LEU
8	H	127	LEU
8	H	196	ARG
9	I	37	ASN
9	I	171	LEU
10	J	2	ASP
10	J	3	ILE
10	J	35	THR
10	J	90	LYS
10	J	99	GLN
11	K	4	LEU
11	K	9	GLN
11	K	104	TYR
11	K	107	LYS
11	K	116	ASP
11	K	140	LEU
11	K	148	LEU
12	L	3	ASN
12	L	23	LEU
12	L	31	THR
12	L	130	SER
12	L	150	LEU
13	M	2	GLN

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Mol	Chain	Res	Type
13	M	48	ASN
13	M	70	LEU
13	M	104	ARG
13	M	161	ARG
13	M	187	ARG
14	N	9	LYS
14	N	83	LYS
14	N	104	ASP
14	N	107	LYS
1	O	62	SER
1	O	122	THR
1	O	157	PHE
1	O	250	LEU
2	P	50	LYS
2	P	55	LEU
2	P	79	LEU
2	P	113	ARG
2	P	191	LEU
2	P	238	LEU
3	Q	4	ARG
3	Q	38	ASN
3	Q	50	LEU
3	Q	51	LYS
3	Q	77	ASN
3	Q	147	GLN
3	Q	160	GLN
3	Q	169	VAL
3	Q	180	LYS
3	Q	206	LYS
3	Q	240	GLU
4	R	99	ILE
4	R	125	LEU
4	R	143	ASP
4	R	176	LEU
4	R	193	LEU
4	R	214	ILE
4	R	235	LEU
4	R	236	LYS
4	R	242	GLU
5	S	9	THR
5	S	29	LYS
5	S	71	LEU

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Mol	Chain	Res	Type
5	S	184	ASN
5	S	188	LEU
5	S	202	ASP
5	S	208	ASP
5	S	231	LYS
6	T	117	GLN
6	T	123	ASN
6	T	139	LYS
6	T	172	LEU
6	T	181	GLU
6	T	201	GLU
6	T	202	ASP
6	T	207	ASP
6	T	214	TRP
6	T	240	GLN
7	U	26	THR
7	U	75	ASN
7	U	83	ASN
7	U	115	LEU
7	U	117	GLN
7	U	122	ARG
7	U	125	MET
7	U	208	GLU
7	U	235	ARG
7	U	236	LEU
8	V	22	GLN
8	V	30	ASN
8	V	56	THR
8	V	68	LEU
8	V	80	LEU
8	V	127	LEU
8	V	196	ARG
9	W	37	ASN
9	W	171	LEU
10	X	2	ASP
10	X	3	ILE
10	X	35	THR
10	X	90	LYS
10	X	99	GLN
11	Y	4	LEU
11	Y	9	GLN
11	Y	104	TYR

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Mol	Chain	Res	Type
11	Y	107	LYS
11	Y	116	ASP
11	Y	140	LEU
11	Y	148	LEU
12	Z	3	ASN
12	Z	23	LEU
12	Z	31	THR
12	Z	130	SER
12	Z	136	CYS
12	Z	150	LEU
13	a	2	GLN
13	a	48	ASN
13	a	70	LEU
13	a	104	ARG
13	a	161	ARG
13	a	187	ARG
14	b	9	LYS
14	b	83	LYS
14	b	104	ASP
14	b	107	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (105) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	94	HIS
2	B	20	GLN
2	B	58	GLN
2	B	95	GLN
2	B	119	GLN
2	B	123	GLN
2	B	155	ASN
3	C	17	GLN
3	C	38	ASN
3	C	77	ASN
3	C	147	GLN
3	C	160	GLN
4	D	15	GLN
4	D	100	ASN
4	D	146	GLN
4	D	225	ASN
5	E	68	HIS
5	E	92	ASN

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Mol	Chain	Res	Type
5	E	99	ASN
5	E	116	GLN
5	E	118	ASN
5	E	120	GLN
5	E	151	ASN
5	E	184	ASN
6	F	19	GLN
6	F	86	ASN
6	F	117	GLN
6	F	123	ASN
6	F	191	GLN
6	F	240	GLN
7	G	83	ASN
7	G	114	ASN
7	G	117	GLN
7	G	121	GLN
7	G	166	GLN
7	G	167	GLN
8	H	22	GLN
8	H	66	HIS
8	H	116	HIS
9	I	37	ASN
10	J	55	GLN
10	J	146	HIS
11	K	85	ASN
11	K	176	ASN
12	L	3	ASN
12	L	49	ASN
12	L	70	ASN
12	L	158	ASN
12	L	159	GLN
13	M	48	ASN
13	M	102	GLN
13	M	194	ASN
13	M	213	GLN
1	O	94	HIS
2	P	20	GLN
2	P	58	GLN
2	P	95	GLN
2	P	119	GLN
2	P	123	GLN
3	Q	17	GLN

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Mol	Chain	Res	Type
3	Q	38	ASN
3	Q	77	ASN
3	Q	147	GLN
3	Q	160	GLN
4	R	15	GLN
4	R	91	HIS
4	R	100	ASN
4	R	146	GLN
4	R	225	ASN
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	151	ASN
5	S	184	ASN
6	T	19	GLN
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	114	ASN
7	U	117	GLN
7	U	121	GLN
7	U	166	GLN
7	U	167	GLN
8	V	22	GLN
8	V	66	HIS
9	W	37	ASN
10	X	55	GLN
10	X	146	HIS
11	Y	85	ASN
11	Y	176	ASN
12	Z	3	ASN
12	Z	49	ASN
12	Z	70	ASN
12	Z	79	HIS
12	Z	158	ASN
13	a	48	ASN

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Mol	Chain	Res	Type
13	a	102	GLN
13	a	194	ASN
13	a	213	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 18 ligands modelled in this entry, 12 are monoatomic - leaving 6 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	4KF	H	301	8	23,23,24	1.82	4 (17%)	25,30,32	1.59	3 (12%)
17	4KF	K	301	11	23,23,24	1.75	2 (8%)	25,30,32	1.39	3 (12%)
17	4KF	N	201	14	23,23,24	1.56	2 (8%)	25,30,32	1.58	5 (20%)
17	4KF	V	301	8	23,23,24	1.73	3 (13%)	25,30,32	1.38	5 (20%)
17	4KF	Y	301	11	23,23,24	1.82	3 (13%)	25,30,32	1.54	6 (24%)
17	4KF	b	201	14	23,23,24	1.84	3 (13%)	25,30,32	1.39	5 (20%)

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	4KF	H	301	8	-	0/25/25/27	0/1/1/1
17	4KF	K	301	11	-	0/25/25/27	0/1/1/1
17	4KF	N	201	14	-	0/25/25/27	0/1/1/1
17	4KF	V	301	8	-	0/25/25/27	0/1/1/1
17	4KF	Y	301	11	-	0/25/25/27	0/1/1/1
17	4KF	b	201	14	-	0/25/25/27	0/1/1/1

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
17	b	201	4KF	C6-C5	-6.87	1.35	1.51
17	Y	301	4KF	C6-C5	-6.79	1.35	1.51
17	K	301	4KF	C6-C5	-6.13	1.37	1.51
17	H	301	4KF	C6-C5	-5.99	1.37	1.51
17	N	201	4KF	C6-C5	-5.75	1.38	1.51
17	V	301	4KF	C6-C5	-5.48	1.38	1.51
17	H	301	4KF	C10-C13	-3.33	1.49	1.55
17	H	301	4KF	C10-C9	-3.26	1.51	1.54
17	K	301	4KF	C10-C13	-3.06	1.50	1.55
17	V	301	4KF	C10-C13	-2.91	1.50	1.55
17	b	201	4KF	C10-C9	-2.78	1.51	1.54
17	b	201	4KF	C10-C13	-2.64	1.51	1.55
17	V	301	4KF	C10-C9	-2.63	1.52	1.54
17	N	201	4KF	C10-C13	-2.51	1.51	1.55
17	H	301	4KF	O21-C20	-2.06	1.33	1.37
17	Y	301	4KF	C10-C13	-2.04	1.52	1.55
17	Y	301	4KF	O2-C1	-2.04	1.33	1.37

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
17	H	301	4KF	C22-O21-C20	-5.36	104.96	117.51
17	N	201	4KF	O17-C9-C8	-3.49	102.36	110.47
17	Y	301	4KF	C14-C13-C15	-3.20	103.73	111.87
17	N	201	4KF	C6-C5-C19	-3.18	114.99	120.62
17	Y	301	4KF	C6-C5-C4	-3.13	115.08	120.62
17	Y	301	4KF	C3-O2-C1	-2.99	110.52	117.51
17	K	301	4KF	C14-C13-C15	-2.91	104.46	111.87
17	H	301	4KF	C6-C5-C19	-2.90	115.48	120.62
17	K	301	4KF	C6-C5-C4	-2.87	115.53	120.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
17	N	201	4KF	C22-O21-C20	-2.81	110.93	117.51
17	V	301	4KF	C6-C5-C19	-2.56	116.08	120.62
17	V	301	4KF	C22-O21-C20	-2.50	111.66	117.51
17	b	201	4KF	C6-C5-C19	-2.42	116.33	120.62
17	b	201	4KF	O12-C11-C10	-2.15	119.17	125.92
17	b	201	4KF	C22-O21-C20	-2.14	112.51	117.51
17	b	201	4KF	C23-C20-C19	-2.06	117.44	120.99
17	Y	301	4KF	O12-C11-C10	-2.05	119.50	125.92
17	V	301	4KF	C5-C6-N7	2.06	117.58	112.88
17	b	201	4KF	C20-C19-C5	2.07	121.88	119.69
17	Y	301	4KF	C5-C6-N7	2.25	118.02	112.88
17	V	301	4KF	C3-O2-C1	2.29	122.87	117.51
17	N	201	4KF	C20-C19-C5	2.35	122.18	119.69
17	K	301	4KF	C6-C5-C19	2.41	124.88	120.62
17	Y	301	4KF	C6-C5-C19	2.55	125.14	120.62
17	N	201	4KF	C6-C5-C4	2.60	125.23	120.62
17	V	301	4KF	C6-C5-C4	2.64	125.30	120.62
17	H	301	4KF	C6-C5-C4	2.87	125.70	120.62

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
17	H	301	4KF	2	0
17	K	301	4KF	7	0
17	Y	301	4KF	5	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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, 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%)	-0.52	3 (1%) 81 55	36, 49, 89, 130	0
1	O	250/250 (100%)	-0.49	4 (1%) 74 47	40, 56, 100, 131	0
2	B	244/258 (94%)	-0.50	5 (2%) 68 39	34, 55, 98, 155	0
2	P	244/258 (94%)	-0.46	5 (2%) 68 39	42, 60, 100, 146	0
3	C	240/254 (94%)	-0.38	7 (2%) 55 26	35, 61, 122, 149	0
3	Q	240/254 (94%)	-0.20	11 (4%) 36 14	47, 74, 157, 183	0
4	D	235/260 (90%)	-0.55	1 (0%) 93 80	39, 60, 95, 138	0
4	R	235/260 (90%)	-0.41	3 (1%) 79 53	48, 70, 109, 143	0
5	E	231/234 (98%)	-0.40	1 (0%) 93 80	44, 69, 111, 153	0
5	S	231/234 (98%)	-0.20	8 (3%) 48 21	45, 77, 134, 176	0
6	F	243/288 (84%)	-0.53	3 (1%) 81 55	45, 64, 116, 140	0
6	T	243/288 (84%)	-0.43	1 (0%) 93 80	46, 72, 133, 170	0
7	G	241/252 (95%)	-0.60	0 100 100	35, 55, 89, 138	0
7	U	241/252 (95%)	-0.56	3 (1%) 81 55	39, 56, 89, 136	0
8	H	226/232 (97%)	-0.50	7 (3%) 52 24	32, 51, 82, 150	0
8	V	226/232 (97%)	-0.50	6 (2%) 58 28	34, 50, 78, 150	0
9	I	204/205 (99%)	-0.75	0 100 100	31, 46, 73, 104	0
9	W	204/205 (99%)	-0.72	1 (0%) 91 76	31, 46, 75, 104	0
10	J	195/198 (98%)	-0.65	3 (1%) 76 49	33, 49, 73, 124	0
10	X	195/198 (98%)	-0.61	2 (1%) 84 60	35, 52, 76, 140	0
11	K	212/212 (100%)	-0.65	0 100 100	28, 48, 70, 90	0
11	Y	212/212 (100%)	-0.66	0 100 100	35, 51, 76, 98	0
12	L	222/222 (100%)	-0.68	1 (0%) 91 76	33, 51, 87, 120	0
12	Z	222/222 (100%)	-0.66	0 100 100	33, 49, 84, 119	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
13	M	229/246 (93%)	-0.60	2 (0%) 85 64	33, 56, 85, 100	0
13	a	232/246 (94%)	-0.55	2 (0%) 85 64	29, 56, 84, 102	0
14	N	196/196 (100%)	-0.67	0 100 100	36, 51, 81, 106	0
14	b	196/196 (100%)	-0.62	1 (0%) 91 76	37, 51, 83, 115	0
All	All	6339/6614 (95%)	-0.53	80 (1%) 79 53	28, 56, 106, 183	0

All (80) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
8	V	224	GLN	7.7
3	Q	49	THR	7.3
8	H	224	GLN	5.0
2	P	219	ALA	4.8
3	Q	50	LEU	4.6
8	H	226	GLU	4.4
5	S	202	ASP	4.2
2	P	220	ASN	4.0
2	P	51	VAL	4.0
8	V	226	GLU	3.9
2	B	221	ASP	3.9
10	X	1	MET	3.6
3	C	49	THR	3.6
8	V	222	ASP	3.6
8	V	223	ILE	3.5
9	W	1	SER	3.4
5	E	202	ASP	3.4
2	P	221	ASP	3.3
10	J	1	MET	3.3
14	b	195	GLN	3.2
4	D	1	ASP	3.1
1	A	2	THR	3.0
10	X	194	ASP	2.9
6	F	244	ASN	2.9
2	P	59	ASP	2.9
13	a	1	THR	2.9
1	O	52	SER	2.8
3	C	50	LEU	2.8
7	U	242	GLN	2.8
8	H	223	ILE	2.8
3	Q	205	ALA	2.7
2	B	220	ASN	2.7

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Mol	Chain	Res	Type	RSRZ
6	T	244	ASN	2.7
8	V	221	CYS	2.7
3	Q	239	GLN	2.7
3	Q	236	GLN	2.7
1	O	2	THR	2.7
8	V	225	GLU	2.6
3	Q	206	LYS	2.6
3	Q	48	SER	2.6
4	R	241	ALA	2.6
4	R	1	ASP	2.6
3	C	206	LYS	2.5
5	S	173	ARG	2.5
5	S	51	ASN	2.5
5	S	180	LYS	2.5
3	Q	240	GLU	2.5
1	O	62	SER	2.5
3	Q	187	GLU	2.5
8	H	222	ASP	2.4
5	S	54	GLU	2.4
3	C	240	GLU	2.4
2	B	51	VAL	2.4
4	R	217	GLN	2.4
5	S	204	SER	2.3
6	F	2	THR	2.3
5	S	165	GLN	2.3
13	a	216	ASN	2.3
7	U	222	ASP	2.3
8	H	198	GLU	2.2
3	C	216	ASP	2.2
2	B	203	SER	2.2
3	C	239	GLN	2.2
2	B	59	ASP	2.2
7	U	241	GLU	2.2
8	H	225	GLU	2.1
10	J	193	ASP	2.1
13	M	47	ASP	2.1
12	L	174	TYR	2.1
10	J	194	ASP	2.1
1	A	62	SER	2.1
1	O	53	SER	2.1
3	Q	238	LYS	2.1
13	M	226	LYS	2.0

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Mol	Chain	Res	Type	RSRZ
5	S	57	SER	2.0
6	F	205	GLU	2.0
3	Q	204	GLY	2.0
1	A	1	MET	2.0
8	H	221	CYS	2.0
3	C	236	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, 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	LLDF	B-factors(Å ²)	Q<0.9
15	MG	W	301	1/1	0.97	0.30	5.69	49,49,49,49	0
15	MG	I	301	1/1	0.97	0.24	5.28	52,52,52,52	0
16	CL	N	203	1/1	0.92	0.21	2.79	52,52,52,52	0
17	4KF	K	301	23/24	0.91	0.25	2.32	37,44,50,54	0
17	4KF	H	301	23/24	0.93	0.24	2.25	32,37,40,46	0
17	4KF	Y	301	23/24	0.92	0.26	2.18	39,44,50,50	0
17	4KF	N	201	23/24	0.90	0.24	1.94	33,47,55,56	0
17	4KF	b	201	23/24	0.89	0.23	1.86	37,46,54,55	0
16	CL	b	202	1/1	0.86	0.20	1.25	52,52,52,52	0
17	4KF	V	301	23/24	0.94	0.20	1.15	35,38,43,48	0
15	MG	Z	301	1/1	0.95	0.14	-0.63	45,45,45,45	0
15	MG	N	202	1/1	0.98	0.10	-0.70	35,35,35,35	0
15	MG	G	301	1/1	0.94	0.10	-0.96	59,59,59,59	0
15	MG	K	302	1/1	0.89	0.09	-1.06	41,41,41,41	0
15	MG	I	302	1/1	0.95	0.07	-1.32	45,45,45,45	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
15	MG	L	301	1/1	0.99	0.07	-2.38	50,50,50,50	0
16	CL	U	301	1/1	0.96	0.20	-	47,47,47,47	0
16	CL	G	302	1/1	0.97	0.28	-	44,44,44,44	0

6.5 Other polymers [i](#)

There are no such residues in this entry.