



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 07:40 am GMT

PDB ID : 4NNN
Title : yCP in complex with MG132
Authors : Stein, M.L.; Cui, H.; Beck, P.; Dubiella, C.; Voss, C.; Krueger, A.; Schmidt, B.; Groll, M.
Deposited on : 2013-11-18
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

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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : recalc28949

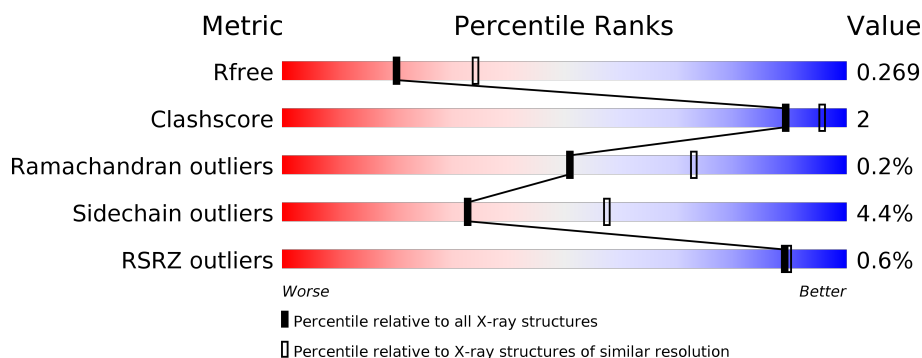
1 Overall quality at a glance

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.











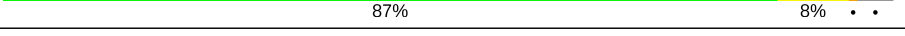












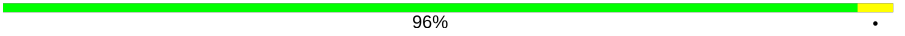
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	100719	3846 (2.50-2.50)
Clashscore	112137	4554 (2.50-2.50)
Ramachandran outliers	110173	4463 (2.50-2.50)
Sidechain outliers	110143	4465 (2.50-2.50)
RSRZ outliers	101464	3876 (2.50-2.50)

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 ≥ 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></div> <div>94% 6%</div> </div>
1	O	250	<div> <div></div> <div>93% 7%</div> </div>
2	B	258	<div> <div></div> <div>84% 10% 5%</div> </div>
2	P	258	<div> <div></div> <div>83% 11% 5%</div> </div>
3	C	254	<div> <div></div> <div>87% 6% • 6%</div> </div>
3	Q	254	<div> <div></div> <div>87% 6% • 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	K	302	-	-	-	X
15	MG	V	302	-	-	-	X
16	ALD	H	301	-	-	-	X
16	ALD	K	301	-	-	-	X
16	ALD	N	201	-	-	-	X
16	ALD	V	301	-	-	-	X
16	ALD	Y	301	-	-	-	X
16	ALD	b	201	-	-	-	X

2 Entry composition

There are 17 unique types of molecules in this entry. The entry contains 50433 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	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	280	312	7			
11	Y	212	Total	C	N	O	S	0	0	0
			1644	1045	280	312	7			

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

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

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

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

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

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

- Molecule 15 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

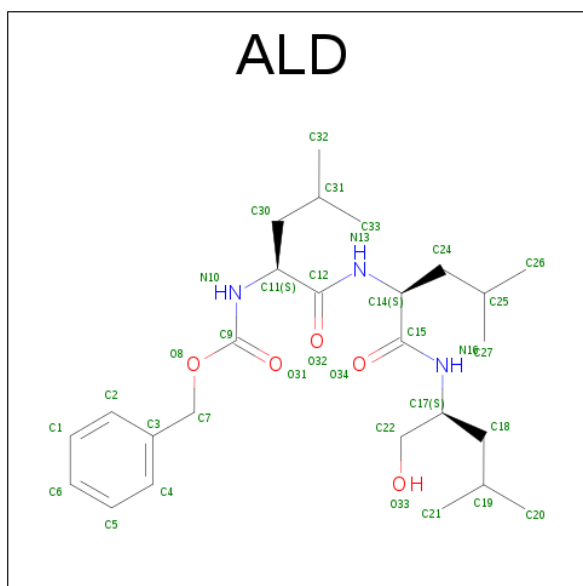
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
15	G	1	Total	Mg	0	0
			1	1		
15	K	1	Total	Mg	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
15	H	1	Total	Mg	0	0
			1	1		
15	I	1	Total	Mg	0	0
			1	1		
15	V	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	Y	1	Total	Mg	0	0
			1	1		

- Molecule 16 is N-[(BENZYLOXY)CARBONYL]-L-LEUCYL-N-[(2S)-1-HYDROXY-4-METHYLPENTAN-2-YL]-L-LEUCINAMIDE (three-letter code: ALD) (formula: C₂₆H₄₃N₃O₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
16	H	1	Total	C	N	O	0	0
			34	26	3	5		
16	K	1	Total	C	N	O	0	0
			34	26	3	5		
16	N	1	Total	C	N	O	0	0
			34	26	3	5		
16	V	1	Total	C	N	O	0	0
			34	26	3	5		
16	Y	1	Total	C	N	O	0	0
			34	26	3	5		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
16	b	1	Total	C	N	O	0	0
			34	26	3	5		

- Molecule 17 is water.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
17	A	36	Total	O			0	0
			36	36				
17	B	31	Total	O			0	0
			31	31				
17	C	27	Total	O			0	0
			27	27				
17	D	22	Total	O			0	0
			22	22				
17	E	11	Total	O			0	0
			11	11				
17	F	32	Total	O			0	0
			32	32				
17	G	50	Total	O			0	0
			50	50				
17	H	31	Total	O			0	0
			31	31				
17	I	46	Total	O			0	0
			46	46				
17	J	33	Total	O			0	0
			33	33				
17	K	40	Total	O			0	0
			40	40				
17	L	41	Total	O			0	0
			41	41				
17	M	41	Total	O			0	0
			41	41				
17	N	43	Total	O			0	0
			43	43				
17	O	24	Total	O			0	0
			24	24				
17	P	22	Total	O			0	0
			22	22				
17	Q	20	Total	O			0	0
			20	20				
17	R	26	Total	O			0	0
			26	26				

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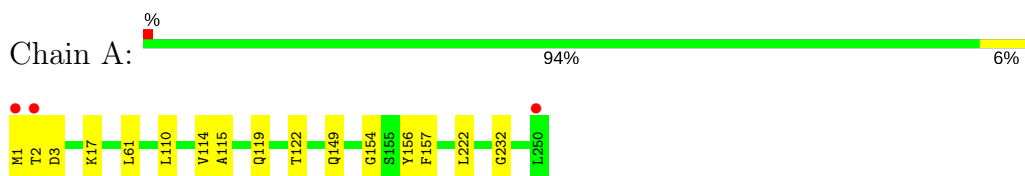
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
17	S	11	Total 11	O 11	0	0
17	T	28	Total 28	O 28	0	0
17	U	45	Total 45	O 45	0	0
17	V	41	Total 41	O 41	0	0
17	W	42	Total 42	O 42	0	0
17	X	31	Total 31	O 31	0	0
17	Y	35	Total 35	O 35	0	0
17	Z	32	Total 32	O 32	0	0
17	a	55	Total 55	O 55	0	0
17	b	29	Total 29	O 29	0	0

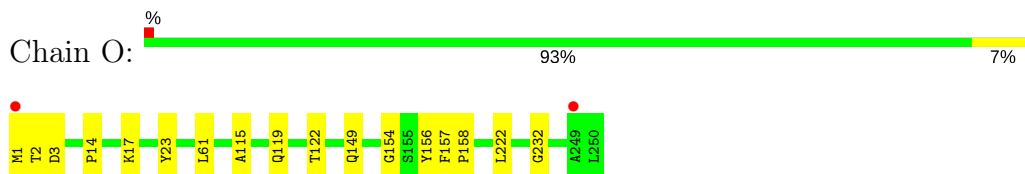
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

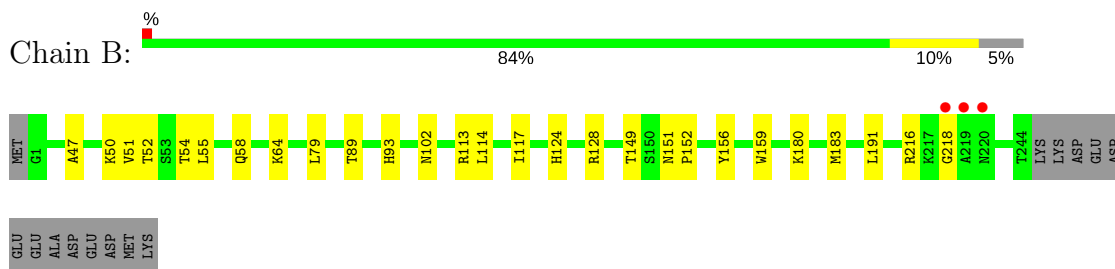
- Molecule 1: Proteasome subunit alpha type-2



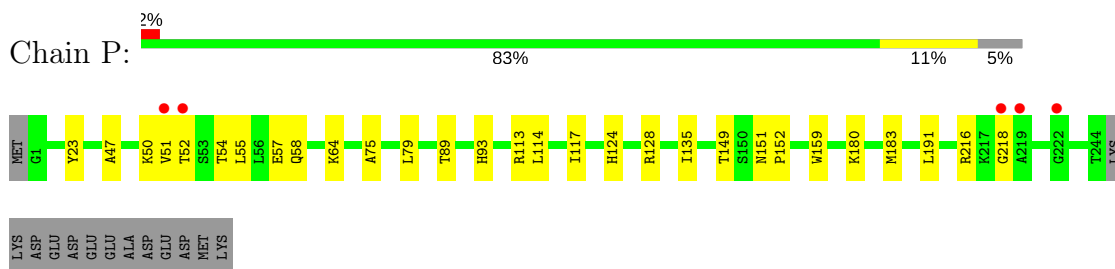
- Molecule 1: Proteasome subunit alpha type-2



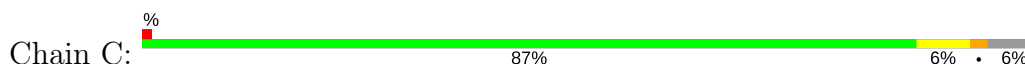
- Molecule 2: Proteasome subunit alpha type-3



- Molecule 2: Proteasome subunit alpha type-3

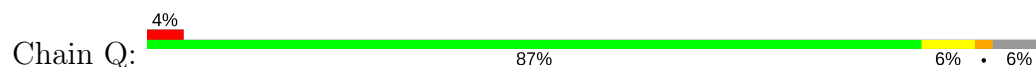


- Molecule 3: Proteasome subunit alpha type-4

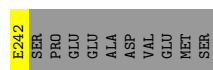
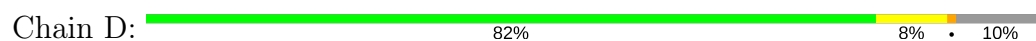




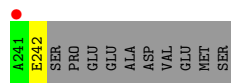
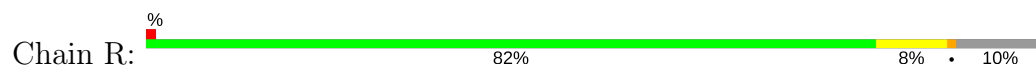
• Molecule 3: Proteasome subunit alpha type-4



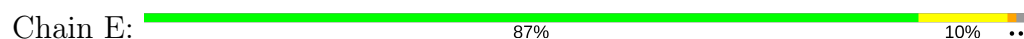
• Molecule 4: Proteasome subunit alpha type-5



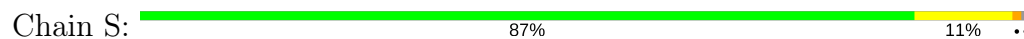
• Molecule 4: Proteasome subunit alpha type-5



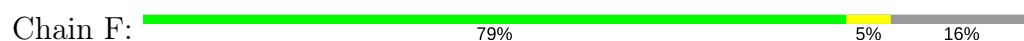
• Molecule 5: Proteasome subunit alpha type-6



• Molecule 5: Proteasome subunit alpha type-6

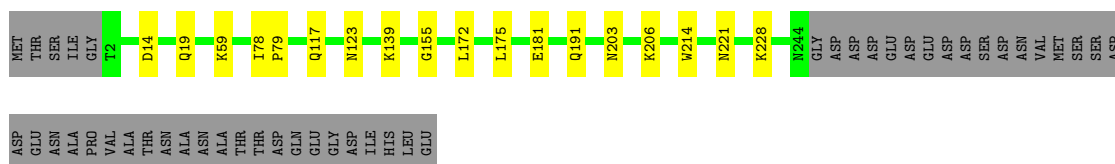


• Molecule 6: Probable proteasome subunit alpha type-7



- Molecule 6: Probable proteasome subunit alpha type-7

Chain T: 78% 6% 16%

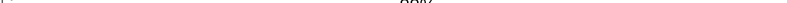


- Molecule 7: Proteasome subunit alpha type-1

Chain G: 88% 7% .

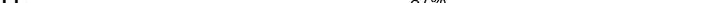


- Molecule 7: Proteasome subunit alpha type-1

Chain U:  88% 8% .



- Molecule 8: Proteasome subunit beta type-2

Chain H:  87% 8% ..



- Molecule 8: Proteasome subunit beta type-2

Chain V: 88% 7% . .

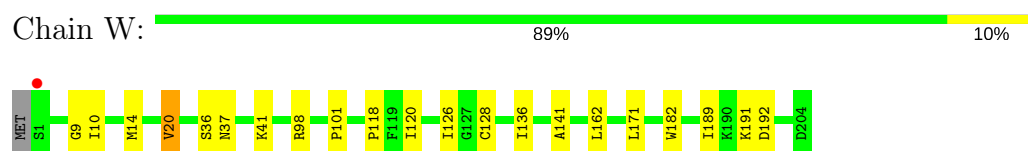


- Molecule 9: Proteasome subunit beta type-3

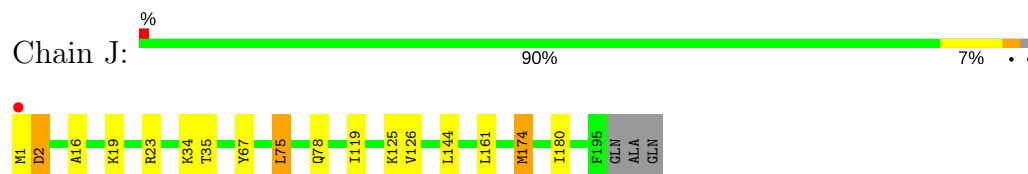
Chain I: 89% 10%



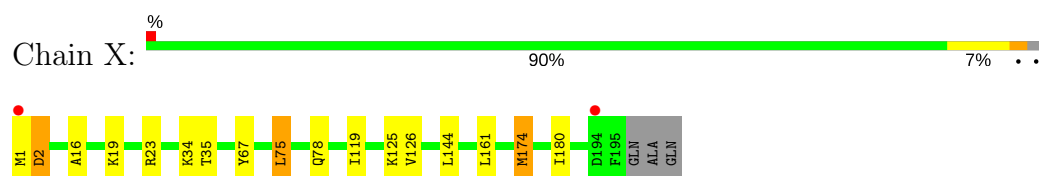
- Molecule 9: Proteasome subunit beta type-3



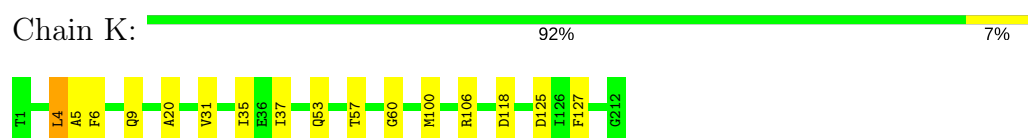
- Molecule 10: Proteasome subunit beta type-4



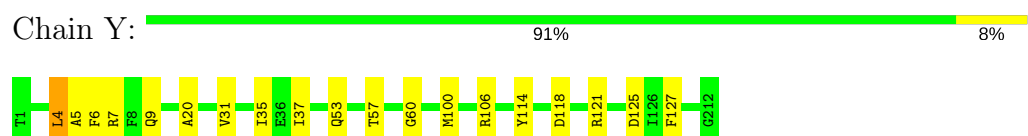
- Molecule 10: Proteasome subunit beta type-4



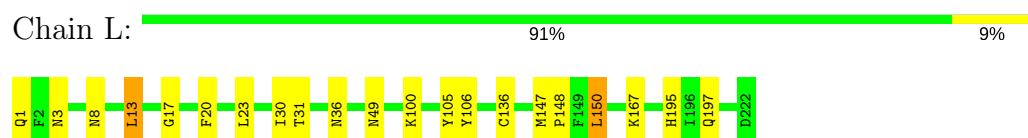
- Molecule 11: Proteasome subunit beta type-5



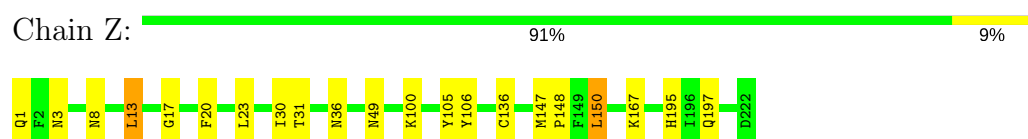
- Molecule 11: Proteasome subunit beta type-5



- Molecule 12: Proteasome subunit beta type-6



- Molecule 12: Proteasome subunit beta type-6



- Molecule 13: Proteasome subunit beta type-7

Chain M:

87%

7%

5%



• Molecule 13: Proteasome subunit beta type-7

Chain a:

91%

•

5%



• Molecule 14: Proteasome subunit beta type-1

Chain N:

%

93%

6%

•

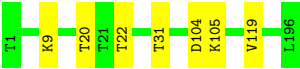


• Molecule 14: Proteasome subunit beta type-1

Chain b:

96%

•



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	134.69Å 299.69Å 144.45Å 90.00° 112.88° 90.00°	Depositor
Resolution (Å)	15.00 – 2.50 15.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	93.9 (15.00-2.50) 93.9 (15.00-2.50)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.75 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.238 , 0.249 0.249 , 0.269	Depositor DCC
R_{free} test set	16930 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	28.4	Xtriage
Anisotropy	1.410	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 15.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	50433	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.29% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.29	0/1952	0.51	0/2642
1	O	0.29	0/1952	0.50	0/2642
2	B	0.29	0/1934	0.52	0/2618
2	P	0.29	0/1934	0.53	0/2618
3	C	0.29	0/1910	0.53	0/2586
3	Q	0.28	0/1910	0.53	0/2586
4	D	0.28	0/1837	0.51	0/2475
4	R	0.28	0/1837	0.51	0/2475
5	E	0.28	0/1800	0.51	0/2433
5	S	0.28	0/1800	0.51	0/2433
6	F	0.29	0/1932	0.49	0/2609
6	T	0.29	0/1932	0.48	0/2609
7	G	0.28	0/1945	0.50	0/2634
7	U	0.29	0/1945	0.51	0/2634
8	H	0.25	0/1715	0.51	0/2326
8	V	0.26	0/1715	0.51	0/2326
9	I	0.28	0/1611	0.51	0/2174
9	W	0.28	0/1611	0.51	0/2174
10	J	0.28	0/1589	0.51	0/2142
10	X	0.29	0/1589	0.51	0/2142
11	K	0.28	0/1681	0.51	1/2274 (0.0%)
11	Y	0.28	0/1681	0.52	1/2274 (0.0%)
12	L	0.28	0/1795	0.50	0/2420
12	Z	0.28	0/1795	0.50	0/2420
13	M	0.28	0/1855	0.53	0/2514
13	a	0.28	0/1855	0.53	0/2514
14	N	0.27	0/1541	0.49	0/2087
14	b	0.27	0/1541	0.50	0/2087
All	All	0.28	0/50194	0.51	2/67868 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	K	4	LEU	CA-CB-CG	5.07	126.95	115.30
11	Y	4	LEU	CA-CB-CG	5.04	126.89	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1915	0	1929	7	0
1	O	1915	0	1929	10	0
2	B	1904	0	1904	12	0
2	P	1904	0	1904	14	0
3	C	1881	0	1895	10	0
3	Q	1881	0	1895	10	0
4	D	1813	0	1797	8	0
4	R	1813	0	1797	8	0
5	E	1773	0	1775	8	0
5	S	1773	0	1775	8	0
6	F	1892	0	1883	4	0
6	T	1892	0	1883	5	0
7	G	1907	0	1901	6	0
7	U	1907	0	1901	9	0
8	H	1684	0	1687	8	0
8	V	1684	0	1687	7	0
9	I	1581	0	1574	11	0
9	W	1581	0	1574	10	0
10	J	1561	0	1569	10	0
10	X	1561	0	1569	10	0
11	K	1644	0	1594	8	0
11	Y	1644	0	1594	10	0
12	L	1757	0	1711	7	0
12	Z	1757	0	1711	7	0
13	M	1824	0	1832	5	0
13	a	1824	0	1832	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
14	N	1512	0	1480	8	0
14	b	1512	0	1480	0	0
15	G	1	0	0	0	0
15	H	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	H	34	0	42	0	0
16	K	34	0	42	0	0
16	N	34	0	42	0	0
16	V	34	0	42	0	0
16	Y	34	0	42	0	0
16	b	34	0	42	0	0
17	A	36	0	0	0	0
17	B	31	0	0	0	0
17	C	27	0	0	0	0
17	D	22	0	0	0	0
17	E	11	0	0	0	0
17	F	32	0	0	0	0
17	G	50	0	0	0	0
17	H	31	0	0	0	0
17	I	46	0	0	0	0
17	J	33	0	0	0	0
17	K	40	0	0	0	0
17	L	41	0	0	0	0
17	M	41	0	0	0	0
17	N	43	0	0	1	0
17	O	24	0	0	0	0
17	P	22	0	0	0	0
17	Q	20	0	0	0	0
17	R	26	0	0	0	0
17	S	11	0	0	0	0
17	T	28	0	0	1	0
17	U	45	0	0	0	0
17	V	41	0	0	0	0
17	W	42	0	0	0	0
17	X	31	0	0	0	0
17	Y	35	0	0	0	0
17	Z	32	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
17	a	55	0	0	0	0
17	b	29	0	0	0	0
All	All	50433	0	49314	201	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 (201) 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)
4:R:99:ILE:HD11	4:R:104:LEU:HB2	1.66	0.77
4:D:99:ILE:HD11	4:D:104:LEU:HB2	1.66	0.75
12:Z:13:LEU:HD13	12:Z:150:LEU:HD21	1.69	0.75
12:L:13:LEU:HD13	12:L:150:LEU:HD21	1.68	0.74
11:K:100:MET:HE3	11:K:127:PHE:HB2	1.72	0.71
10:X:1:MET:O	10:X:2:ASP:HB2	1.89	0.70
10:J:1:MET:O	10:J:2:ASP:HB2	1.93	0.67
11:Y:100:MET:HE3	11:Y:127:PHE:HB2	1.75	0.67
7:U:23:PHE:O	7:U:26:THR:HB	1.96	0.66
11:Y:100:MET:CE	11:Y:127:PHE:HB2	2.26	0.65
7:G:23:PHE:O	7:G:26:THR:HB	1.98	0.64
11:K:100:MET:CE	11:K:127:PHE:HB2	2.26	0.64
3:Q:51:LYS:O	3:Q:52:LEU:HB2	1.98	0.64
11:Y:5:ALA:HB3	11:Y:100:MET:HE2	1.81	0.63
3:C:51:LYS:O	3:C:52:LEU:HB2	1.99	0.62
9:I:98:ARG:O	9:I:126:ILE:HD11	1.99	0.62
9:W:98:ARG:O	9:W:126:ILE:HD11	2.01	0.61
1:A:122:THR:HG22	2:B:128:ARG:HH21	1.66	0.60
14:N:20:THR:CG2	14:N:28:ASN:HB3	2.31	0.60
11:K:5:ALA:HB3	11:K:100:MET:HE2	1.84	0.60
14:N:152:VAL:HA	14:N:175:MET:HE1	1.83	0.60
10:J:16:ALA:HB2	10:J:161:LEU:HD21	1.85	0.59
1:O:122:THR:HG22	2:P:128:ARG:HH21	1.67	0.58
2:P:93:HIS:HB3	2:P:113:ARG:HH21	1.69	0.57
2:B:93:HIS:HB3	2:B:113:ARG:HH21	1.69	0.57
10:J:1:MET:HA	10:J:34:LYS:HE3	1.86	0.57
10:X:16:ALA:HB2	10:X:161:LEU:HD21	1.87	0.57
10:X:1:MET:HA	10:X:34:LYS:HE3	1.87	0.56
9:I:20:VAL:HG13	9:I:118:PRO:HB3	1.88	0.56
3:C:160:GLN:HE21	3:C:160:GLN:HA	1.71	0.55
1:O:119:GLN:O	1:O:122:THR:HB	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:124:HIS:HB3	3:C:124:VAL:HG12	1.88	0.55
5:S:9:THR:HG21	5:S:119:THR:HA	1.89	0.55
8:H:104:ASP:HB2	8:H:105:PRO:HD2	1.89	0.55
2:P:124:HIS:HB3	3:Q:124:VAL:HG12	1.88	0.55
3:Q:160:GLN:HE21	3:Q:160:GLN:HA	1.71	0.55
1:A:149:GLN:O	1:A:156:TYR:HA	2.07	0.55
5:E:9:THR:HG21	5:E:119:THR:HA	1.89	0.55
4:R:88:ALA:HA	4:R:99:ILE:HG21	1.89	0.55
11:Y:53:GLN:O	11:Y:57:THR:HG23	2.07	0.54
4:D:88:ALA:HA	4:D:99:ILE:HG21	1.89	0.54
5:E:12:PHE:H	6:F:19:GLN:HE22	1.54	0.54
8:V:104:ASP:HB2	8:V:105:PRO:HD2	1.89	0.54
10:X:1:MET:O	10:X:2:ASP:CB	2.56	0.54
9:W:20:VAL:HG13	9:W:118:PRO:HB3	1.90	0.53
1:A:119:GLN:O	1:A:122:THR:HB	2.08	0.53
1:O:149:GLN:O	1:O:156:TYR:HA	2.08	0.53
11:K:53:GLN:O	11:K:57:THR:HG23	2.09	0.52
3:C:201:VAL:O	3:C:202:GLN:CB	2.58	0.52
8:V:52:THR:O	8:V:56:THR:HB	2.09	0.52
8:H:52:THR:O	8:H:56:THR:HB	2.10	0.51
9:W:9:GLY:HA3	9:W:41:LYS:HE2	1.93	0.51
9:I:9:GLY:HA3	9:I:41:LYS:HE2	1.91	0.51
10:J:1:MET:O	10:J:2:ASP:CB	2.57	0.51
9:W:10:ILE:HG21	9:W:141:ALA:HB3	1.93	0.51
9:I:36:SER:HB2	10:J:126:VAL:HG11	1.92	0.51
3:Q:9:PHE:H	4:R:15:GLN:HE22	1.57	0.51
5:S:12:PHE:H	6:T:19:GLN:HE22	1.58	0.51
3:Q:201:VAL:O	3:Q:202:GLN:CB	2.59	0.50
9:I:10:ILE:HG21	9:I:141:ALA:HB3	1.93	0.50
7:U:187:GLU:HG2	7:U:192:LYS:HB2	1.92	0.50
12:L:31:THR:HG23	12:L:36:ASN:HD21	1.76	0.50
2:B:151:ASN:HB2	2:B:152:PRO:HD2	1.93	0.50
3:C:201:VAL:HG13	3:C:202:GLN:N	2.27	0.50
2:P:151:ASN:HB2	2:P:152:PRO:HD2	1.94	0.49
9:W:36:SER:HB2	10:X:126:VAL:HG11	1.93	0.49
10:J:19:LYS:HD3	10:J:180:ILE:HG13	1.93	0.49
5:E:68:HIS:HE1	5:E:102:LEU:O	1.94	0.49
7:G:187:GLU:HG2	7:G:192:LYS:HB2	1.92	0.49
10:J:119:ILE:HG12	10:J:125:LYS:HG3	1.94	0.49
5:S:68:HIS:HE1	5:S:102:LEU:O	1.94	0.49
10:J:174:MET:HA	10:X:174:MET:HA	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:Q:198:LEU:HA	3:Q:201:VAL:HG12	1.94	0.49
3:C:198:LEU:HA	3:C:201:VAL:HG12	1.95	0.49
3:Q:201:VAL:HG13	3:Q:202:GLN:N	2.28	0.49
10:X:119:ILE:HG12	10:X:125:LYS:HG3	1.95	0.49
7:U:195:GLU:HG3	7:U:235:ARG:HG3	1.94	0.48
1:O:23:TYR:CD1	7:U:12:PRO:HA	2.48	0.48
10:X:19:LYS:HD3	10:X:180:ILE:HG13	1.94	0.48
5:S:87:LEU:HD21	5:S:107:ALA:HB1	1.96	0.48
8:V:50:ALA:HB2	9:W:128:CYS:HB2	1.96	0.48
12:Z:17:GLY:HA3	12:Z:20:PHE:CE1	2.49	0.48
5:E:131:LEU:HB2	5:E:146:PHE:HB3	1.94	0.48
3:C:201:VAL:O	3:C:202:GLN:HB2	2.13	0.48
4:R:160:ASN:HB3	4:R:179:TRP:CE2	2.49	0.48
2:B:151:ASN:HB2	2:B:152:PRO:CD	2.43	0.48
5:E:87:LEU:HD21	5:E:107:ALA:HB1	1.95	0.48
8:H:148:LYS:HE3	8:H:177:VAL:HG11	1.95	0.48
3:Q:201:VAL:O	3:Q:202:GLN:HB2	2.14	0.48
7:G:195:GLU:HG3	7:G:235:ARG:HG3	1.95	0.48
5:S:131:LEU:HB2	5:S:146:PHE:HB3	1.95	0.48
12:L:17:GLY:HA3	12:L:20:PHE:CE1	2.48	0.47
13:M:127:LEU:HG	13:M:142:LEU:HD12	1.95	0.47
12:Z:31:THR:HG23	12:Z:36:ASN:HD21	1.78	0.47
2:B:47:ALA:HB1	2:B:64:LYS:HD2	1.97	0.47
11:K:6:PHE:HA	11:K:125:ASP:O	2.14	0.47
2:P:47:ALA:HB1	2:P:64:LYS:HD2	1.97	0.47
9:W:20:VAL:HG23	9:W:189:ILE:HB	1.96	0.47
10:J:1:MET:HA	10:J:34:LYS:CE	2.45	0.47
8:V:148:LYS:HE3	8:V:177:VAL:HG11	1.95	0.47
2:P:151:ASN:HB2	2:P:152:PRO:CD	2.44	0.47
2:P:149:THR:HG1	2:P:159:TRP:HE1	1.63	0.47
2:P:180:LYS:O	2:P:183:MET:HB2	2.14	0.47
5:S:175:LEU:HA	5:S:178:PHE:CE2	2.50	0.47
9:I:20:VAL:HG23	9:I:189:ILE:HB	1.96	0.47
11:K:37:ILE:HG23	11:K:60:GLY:HA2	1.97	0.46
4:D:160:ASN:HB3	4:D:179:TRP:CE2	2.50	0.46
12:L:195:HIS:HD2	12:L:197:GLN:H	1.62	0.46
12:Z:195:HIS:HD2	12:Z:197:GLN:H	1.62	0.46
11:Y:6:PHE:HA	11:Y:125:ASP:O	2.15	0.46
12:Z:147:MET:N	12:Z:148:PRO:HD2	2.31	0.46
3:C:9:PHE:H	4:D:15:GLN:HE22	1.62	0.46
12:L:8:ASN:HA	12:L:30:ILE:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:M:27:LEU:HD12	13:M:36:PHE:O	2.15	0.46
5:E:175:LEU:HA	5:E:178:PHE:CE2	2.51	0.46
11:K:5:ALA:HB3	11:K:100:MET:CE	2.45	0.46
12:L:100:LYS:HD3	12:L:105:TYR:CZ	2.51	0.46
11:Y:5:ALA:HB3	11:Y:100:MET:CE	2.44	0.45
9:I:101:PRO:HB3	9:I:126:ILE:HD12	1.99	0.45
10:X:1:MET:HA	10:X:34:LYS:CE	2.45	0.45
6:F:228:LYS:HB2	6:F:228:LYS:HE3	1.68	0.45
8:H:50:ALA:HB2	9:I:128:CYS:HB2	1.99	0.45
9:I:14:MET:HB3	9:I:162:LEU:HD11	1.98	0.45
2:B:149:THR:HG1	2:B:159:TRP:HE1	1.63	0.45
12:Z:100:LYS:HD3	12:Z:105:TYR:CZ	2.52	0.45
6:F:175:LEU:HD21	6:F:191:GLN:NE2	2.32	0.45
12:L:147:MET:N	12:L:148:PRO:HD2	2.31	0.45
2:P:89:THR:HG21	2:P:117:ILE:HD13	1.99	0.45
14:N:67:THR:HA	14:N:71:GLY:O	2.17	0.45
1:O:158:PRO:HB2	2:P:57:GLU:HB3	1.98	0.45
4:D:91:HIS:HB3	4:D:99:ILE:CG2	2.47	0.44
4:R:138:GLY:HA2	4:R:214:ILE:HG12	1.99	0.44
7:U:68:ARG:O	7:U:223:LYS:HA	2.17	0.44
4:D:138:GLY:HA2	4:D:214:ILE:HG12	1.98	0.44
10:J:67:TYR:CE1	10:J:75:LEU:HD13	2.52	0.44
2:B:180:LYS:O	2:B:183:MET:HB2	2.18	0.44
14:N:20:THR:HG22	14:N:31:THR:OG1	2.17	0.44
5:S:42:HIS:HB2	5:S:188:LEU:HD12	1.99	0.44
11:Y:37:ILE:HG23	11:Y:60:GLY:HA2	1.98	0.44
14:N:35:THR:HG22	17:N:305:HOH:O	2.17	0.44
14:N:35:THR:CG2	14:N:45:ARG:HE	2.29	0.44
1:O:222:LEU:HD13	1:O:232:GLY:HA2	1.99	0.44
2:P:89:THR:HG21	2:P:117:ILE:CD1	2.47	0.44
4:R:91:HIS:HB3	4:R:99:ILE:CG2	2.47	0.44
1:A:115:ALA:HB1	1:A:154:GLY:O	2.18	0.44
9:W:101:PRO:HB3	9:W:126:ILE:HD12	2.00	0.43
1:A:222:LEU:HD13	1:A:232:GLY:HA2	2.01	0.43
2:B:89:THR:HG21	2:B:117:ILE:CD1	2.48	0.43
12:Z:8:ASN:HA	12:Z:30:ILE:O	2.18	0.43
7:G:78:ILE:N	7:G:79:PRO:CD	2.81	0.43
6:T:175:LEU:HD21	6:T:191:GLN:NE2	2.33	0.43
7:U:26:THR:HG21	7:U:131:ILE:HD12	2.00	0.43
9:W:14:MET:HB3	9:W:162:LEU:HD11	1.98	0.43
5:E:42:HIS:HB2	5:E:188:LEU:HD12	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:U:78:ILE:N	7:U:79:PRO:CD	2.82	0.43
1:O:1:MET:CG	1:O:2:THR:H	2.32	0.43
10:X:67:TYR:CE1	10:X:75:LEU:HD13	2.54	0.43
3:C:202:GLN:HG3	3:C:203:THR:N	2.34	0.43
9:I:120:ILE:HD12	9:I:136:ILE:HG12	2.00	0.43
14:N:20:THR:HG23	14:N:28:ASN:HB3	1.99	0.43
4:R:73:LEU:HD12	4:R:131:GLY:HA3	2.01	0.43
2:B:216:ARG:HB3	2:B:218:GLY:H	1.84	0.42
2:B:149:THR:O	2:B:156:TYR:HA	2.19	0.42
1:O:115:ALA:HB1	1:O:154:GLY:O	2.19	0.42
8:V:104:ASP:OD1	8:V:106:THR:HB	2.19	0.42
2:B:89:THR:HG21	2:B:117:ILE:HD13	2.00	0.42
8:H:104:ASP:OD1	8:H:106:THR:HB	2.19	0.42
7:G:68:ARG:O	7:G:223:LYS:HA	2.20	0.42
6:T:155:GLY:HA3	7:U:59:THR:HG21	2.02	0.42
11:K:20:ALA:HB2	11:K:31:VAL:HG21	2.02	0.42
1:O:14:PRO:HA	2:P:23:TYR:CD1	2.55	0.42
3:Q:202:GLN:HG3	3:Q:203:THR:N	2.34	0.42
1:A:1:MET:CG	1:A:2:THR:H	2.32	0.42
14:N:35:THR:HG21	14:N:45:ARG:HE	1.85	0.42
11:Y:100:MET:HE1	11:Y:127:PHE:HB2	2.02	0.42
7:G:26:THR:HG21	7:G:131:ILE:HD12	2.01	0.42
1:O:23:TYR:CE1	7:U:12:PRO:HA	2.55	0.42
1:A:110:LEU:O	1:A:114:VAL:HG23	2.19	0.42
9:W:120:ILE:HD12	9:W:136:ILE:HG12	2.01	0.42
11:Y:20:ALA:HB2	11:Y:31:VAL:HG21	2.02	0.41
4:D:30:ILE:HD12	4:D:196:LEU:HG	2.03	0.41
8:H:196:ARG:NH2	9:I:150:GLU:O	2.53	0.41
6:T:19:GLN:NE2	17:T:310:HOH:O	2.52	0.41
8:H:84:LYS:HG3	8:H:85:GLN:N	2.35	0.41
13:M:25:ASP:HA	13:M:195:PHE:CB	2.50	0.41
8:V:3:ILE:HG22	8:V:99:ILE:HD12	2.02	0.41
2:P:216:ARG:HB3	2:P:218:GLY:H	1.85	0.41
13:M:128:ARG:HH11	13:M:138:SER:HB2	1.85	0.41
3:Q:169:VAL:HG23	3:Q:196:SER:HB2	2.02	0.41
4:R:30:ILE:HD12	4:R:196:LEU:HG	2.03	0.41
11:Y:114:TYR:O	11:Y:121:ARG:HA	2.21	0.41
4:D:73:LEU:HD12	4:D:131:GLY:HA3	2.03	0.41
13:M:43:ILE:HA	13:M:44:PRO:HD3	1.95	0.41
5:E:65:CYS:SG	5:E:71:LEU:HD13	2.61	0.41
6:F:78:ILE:HB	6:F:79:PRO:HD3	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:V:112:SER:OG	8:V:120:ASP:HB2	2.21	0.40
3:C:169:VAL:HG23	3:C:196:SER:HB2	2.02	0.40
8:H:112:SER:OG	8:H:120:ASP:HB2	2.21	0.40
5:S:62:ILE:HG21	5:S:213:ALA:HB2	2.04	0.40
2:P:75:ALA:HB3	2:P:135:ILE:HB	2.02	0.40
6:T:78:ILE:HB	6:T:79:PRO:HD3	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	248/250 (99%)	241 (97%)	6 (2%)	1 (0%)	38	59
1	O	248/250 (99%)	241 (97%)	6 (2%)	1 (0%)	38	59
2	B	242/258 (94%)	234 (97%)	7 (3%)	1 (0%)	38	59
2	P	242/258 (94%)	234 (97%)	7 (3%)	1 (0%)	38	59
3	C	238/254 (94%)	226 (95%)	9 (4%)	3 (1%)	14	25
3	Q	238/254 (94%)	227 (95%)	8 (3%)	3 (1%)	14	25
4	D	231/260 (89%)	225 (97%)	6 (3%)	0	100	100
4	R	231/260 (89%)	226 (98%)	5 (2%)	0	100	100
5	E	229/234 (98%)	220 (96%)	9 (4%)	0	100	100
5	S	229/234 (98%)	221 (96%)	8 (4%)	0	100	100
6	F	241/288 (84%)	233 (97%)	8 (3%)	0	100	100
6	T	241/288 (84%)	233 (97%)	8 (3%)	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%)	213 (97%)	7 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
8	V	220/232 (95%)	213 (97%)	7 (3%)	0	100	100
9	I	202/205 (98%)	195 (96%)	7 (4%)	0	100	100
9	W	202/205 (98%)	196 (97%)	6 (3%)	0	100	100
10	J	193/198 (98%)	186 (96%)	6 (3%)	1 (0%)	32	53
10	X	193/198 (98%)	187 (97%)	5 (3%)	1 (0%)	32	53
11	K	210/212 (99%)	205 (98%)	5 (2%)	0	100	100
11	Y	210/212 (99%)	204 (97%)	6 (3%)	0	100	100
12	L	220/222 (99%)	215 (98%)	5 (2%)	0	100	100
12	Z	220/222 (99%)	215 (98%)	5 (2%)	0	100	100
13	M	231/246 (94%)	222 (96%)	9 (4%)	0	100	100
13	a	231/246 (94%)	223 (96%)	8 (4%)	0	100	100
14	N	194/196 (99%)	187 (96%)	7 (4%)	0	100	100
14	b	194/196 (99%)	187 (96%)	7 (4%)	0	100	100
All	All	6276/6614 (95%)	6079 (97%)	185 (3%)	12 (0%)	51	73

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	51	VAL
3	C	202	GLN
10	J	2	ASP
2	P	51	VAL
3	Q	202	GLN
10	X	2	ASP
3	C	205	ALA
3	Q	205	ALA
1	A	3	ASP
3	C	183	PRO
1	O	3	ASP
3	Q	183	PRO

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%)	71	90
1	O	209/209 (100%)	206 (99%)	3 (1%)	71	90
2	B	203/216 (94%)	194 (96%)	9 (4%)	33	57
2	P	203/216 (94%)	195 (96%)	8 (4%)	37	63
3	C	212/226 (94%)	204 (96%)	8 (4%)	38	64
3	Q	212/226 (94%)	204 (96%)	8 (4%)	38	64
4	D	194/215 (90%)	183 (94%)	11 (6%)	24	44
4	R	194/215 (90%)	183 (94%)	11 (6%)	24	44
5	E	190/193 (98%)	175 (92%)	15 (8%)	14	27
5	S	190/193 (98%)	176 (93%)	14 (7%)	16	30
6	F	201/239 (84%)	190 (94%)	11 (6%)	25	46
6	T	201/239 (84%)	189 (94%)	12 (6%)	22	41
7	G	206/210 (98%)	197 (96%)	9 (4%)	33	57
7	U	206/210 (98%)	198 (96%)	8 (4%)	37	63
8	H	181/190 (95%)	172 (95%)	9 (5%)	28	51
8	V	181/190 (95%)	172 (95%)	9 (5%)	28	51
9	I	172/173 (99%)	166 (96%)	6 (4%)	41	68
9	W	172/173 (99%)	166 (96%)	6 (4%)	41	68
10	J	173/175 (99%)	167 (96%)	6 (4%)	41	68
10	X	173/175 (99%)	167 (96%)	6 (4%)	41	68
11	K	169/169 (100%)	164 (97%)	5 (3%)	46	74
11	Y	169/169 (100%)	163 (96%)	6 (4%)	40	67
12	L	185/185 (100%)	176 (95%)	9 (5%)	29	52
12	Z	185/185 (100%)	176 (95%)	9 (5%)	29	52
13	M	199/208 (96%)	190 (96%)	9 (4%)	32	56
13	a	199/208 (96%)	190 (96%)	9 (4%)	32	56
14	N	162/162 (100%)	155 (96%)	7 (4%)	33	58
14	b	162/162 (100%)	155 (96%)	7 (4%)	33	58
All	All	5312/5540 (96%)	5079 (96%)	233 (4%)	33	57

All (233) 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	52	THR
2	B	54	THR
2	B	55	LEU
2	B	58	GLN
2	B	79	LEU
2	B	102	ASN
2	B	114	LEU
2	B	191	LEU
3	C	4	ARG
3	C	38	ASN
3	C	61	LYS
3	C	147	GLN
3	C	160	GLN
3	C	169	VAL
3	C	180	LYS
3	C	203	THR
4	D	20	LEU
4	D	40	LEU
4	D	51	LEU
4	D	99	ILE
4	D	102	GLU
4	D	117	GLU
4	D	176	LEU
4	D	193	LEU
4	D	214	ILE
4	D	236	LYS
4	D	242	GLU
5	E	8	ASP
5	E	9	THR
5	E	10	VAL
5	E	25	LEU
5	E	29	LYS
5	E	54	GLU
5	E	55	LEU
5	E	71	LEU
5	E	99	ASN
5	E	174	THR
5	E	184	ASN
5	E	188	LEU

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Mol	Chain	Res	Type
5	E	207	VAL
5	E	208	ASP
5	E	231	LYS
6	F	59	LYS
6	F	117	GLN
6	F	123	ASN
6	F	139	LYS
6	F	172	LEU
6	F	181	GLU
6	F	203	ASN
6	F	206	LYS
6	F	214	TRP
6	F	221	ASN
6	F	228	LYS
7	G	83	ASN
7	G	115	LEU
7	G	125	MET
7	G	154	TYR
7	G	166	GLN
7	G	181	LYS
7	G	207	THR
7	G	235	ARG
7	G	236	LEU
8	H	3	ILE
8	H	22	GLN
8	H	30	ASN
8	H	34	LEU
8	H	43	CYS
8	H	56	THR
8	H	68	LEU
8	H	127	LEU
8	H	196	ARG
9	I	20	VAL
9	I	37	ASN
9	I	171	LEU
9	I	182	TRP
9	I	191	LYS
9	I	192	ASP
10	J	23	ARG
10	J	35	THR
10	J	75	LEU
10	J	78	GLN

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Mol	Chain	Res	Type
10	J	144	LEU
10	J	174	MET
11	K	4	LEU
11	K	9	GLN
11	K	35	ILE
11	K	106	ARG
11	K	118	ASP
12	L	1	GLN
12	L	3	ASN
12	L	13	LEU
12	L	23	LEU
12	L	49	ASN
12	L	106	TYR
12	L	136	CYS
12	L	150	LEU
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
13	M	204	THR
13	M	212	LEU
13	M	223	LYS
14	N	9	LYS
14	N	20	THR
14	N	22	THR
14	N	31	THR
14	N	104	ASP
14	N	105	LYS
14	N	119	VAL
1	O	17	LYS
1	O	61	LEU
1	O	157	PHE
2	P	50	LYS
2	P	52	THR
2	P	54	THR
2	P	55	LEU
2	P	58	GLN
2	P	79	LEU
2	P	114	LEU

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Mol	Chain	Res	Type
2	P	191	LEU
3	Q	4	ARG
3	Q	38	ASN
3	Q	61	LYS
3	Q	147	GLN
3	Q	160	GLN
3	Q	169	VAL
3	Q	180	LYS
3	Q	203	THR
4	R	20	LEU
4	R	40	LEU
4	R	51	LEU
4	R	99	ILE
4	R	102	GLU
4	R	117	GLU
4	R	176	LEU
4	R	193	LEU
4	R	214	ILE
4	R	236	LYS
4	R	242	GLU
5	S	9	THR
5	S	10	VAL
5	S	25	LEU
5	S	29	LYS
5	S	54	GLU
5	S	55	LEU
5	S	71	LEU
5	S	99	ASN
5	S	174	THR
5	S	184	ASN
5	S	188	LEU
5	S	207	VAL
5	S	208	ASP
5	S	231	LYS
6	T	14	ASP
6	T	59	LYS
6	T	117	GLN
6	T	123	ASN
6	T	139	LYS
6	T	172	LEU
6	T	181	GLU
6	T	203	ASN

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Mol	Chain	Res	Type
6	T	206	LYS
6	T	214	TRP
6	T	221	ASN
6	T	228	LYS
7	U	83	ASN
7	U	115	LEU
7	U	125	MET
7	U	154	TYR
7	U	166	GLN
7	U	181	LYS
7	U	235	ARG
7	U	236	LEU
8	V	3	ILE
8	V	22	GLN
8	V	30	ASN
8	V	34	LEU
8	V	43	CYS
8	V	56	THR
8	V	68	LEU
8	V	127	LEU
8	V	196	ARG
9	W	20	VAL
9	W	37	ASN
9	W	171	LEU
9	W	182	TRP
9	W	191	LYS
9	W	192	ASP
10	X	23	ARG
10	X	35	THR
10	X	75	LEU
10	X	78	GLN
10	X	144	LEU
10	X	174	MET
11	Y	4	LEU
11	Y	7	ARG
11	Y	9	GLN
11	Y	35	ILE
11	Y	106	ARG
11	Y	118	ASP
12	Z	1	GLN
12	Z	3	ASN
12	Z	13	LEU

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Mol	Chain	Res	Type
12	Z	23	LEU
12	Z	49	ASN
12	Z	106	TYR
12	Z	136	CYS
12	Z	150	LEU
12	Z	167	LYS
13	a	43	ILE
13	a	48	ASN
13	a	70	LEU
13	a	104	ARG
13	a	161	ARG
13	a	187	ARG
13	a	204	THR
13	a	212	LEU
13	a	223	LYS
14	b	9	LYS
14	b	20	THR
14	b	22	THR
14	b	31	THR
14	b	104	ASP
14	b	105	LYS
14	b	119	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (120) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	94	HIS
2	B	20	GLN
2	B	95	GLN
2	B	119	GLN
2	B	123	GLN
2	B	176	GLN
3	C	38	ASN
3	C	77	ASN
3	C	116	GLN
3	C	120	GLN
3	C	147	GLN
3	C	160	GLN
4	D	15	GLN
4	D	100	ASN
4	D	146	GLN
4	D	198	GLN

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Mol	Chain	Res	Type
4	D	210	GLN
4	D	225	ASN
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	147	GLN
5	E	184	ASN
6	F	19	GLN
6	F	86	ASN
6	F	117	GLN
6	F	191	GLN
6	F	240	GLN
7	G	6	HIS
7	G	30	ASN
7	G	83	ASN
7	G	114	ASN
7	G	117	GLN
7	G	121	GLN
7	G	167	GLN
8	H	30	ASN
8	H	57	GLN
8	H	66	HIS
8	H	165	ASN
8	H	172	ASN
8	H	189	ASN
10	J	55	GLN
10	J	118	GLN
11	K	85	ASN
11	K	176	ASN
12	L	1	GLN
12	L	3	ASN
12	L	49	ASN
12	L	70	ASN
12	L	165	ASN
13	M	2	GLN
13	M	18	ASN
13	M	48	ASN
13	M	108	ASN
13	M	179	ASN

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Mol	Chain	Res	Type
13	M	194	ASN
13	M	213	GLN
14	N	161	GLN
1	O	94	HIS
2	P	20	GLN
2	P	95	GLN
2	P	119	GLN
2	P	123	GLN
2	P	167	ASN
2	P	176	GLN
3	Q	38	ASN
3	Q	77	ASN
3	Q	147	GLN
3	Q	160	GLN
4	R	15	GLN
4	R	100	ASN
4	R	146	GLN
4	R	198	GLN
4	R	225	ASN
5	S	68	HIS
5	S	92	ASN
5	S	99	ASN
5	S	116	GLN
5	S	120	GLN
5	S	147	GLN
5	S	184	ASN
6	T	19	GLN
6	T	86	ASN
6	T	117	GLN
6	T	191	GLN
6	T	240	GLN
7	U	30	ASN
7	U	83	ASN
7	U	114	ASN
7	U	117	GLN
7	U	121	GLN
8	V	30	ASN
8	V	57	GLN
8	V	66	HIS
8	V	165	ASN
8	V	172	ASN
8	V	189	ASN

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Mol	Chain	Res	Type
10	X	55	GLN
10	X	86	GLN
10	X	118	GLN
11	Y	9	GLN
11	Y	85	ASN
11	Y	176	ASN
12	Z	1	GLN
12	Z	3	ASN
12	Z	49	ASN
12	Z	70	ASN
12	Z	159	GLN
12	Z	165	ASN
13	a	2	GLN
13	a	18	ASN
13	a	48	ASN
13	a	108	ASN
13	a	179	ASN
13	a	194	ASN
13	a	213	GLN
14	b	161	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 14 ligands modelled in this entry, 8 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
16	ALD	H	301	8	34,34,34	0.96	1 (2%)	44,44,44	1.09	3 (6%)
16	ALD	K	301	11	34,34,34	0.93	1 (2%)	44,44,44	1.15	4 (9%)
16	ALD	N	201	14	34,34,34	0.88	1 (2%)	44,44,44	1.23	3 (6%)
16	ALD	V	301	8	34,34,34	0.95	1 (2%)	44,44,44	1.11	3 (6%)
16	ALD	Y	301	11	34,34,34	0.92	1 (2%)	44,44,44	1.22	3 (6%)
16	ALD	b	201	14	34,34,34	0.93	1 (2%)	44,44,44	1.15	3 (6%)

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
16	ALD	H	301	8	-	0/39/39/39	0/1/1/1
16	ALD	K	301	11	-	0/39/39/39	0/1/1/1
16	ALD	N	201	14	-	0/39/39/39	0/1/1/1
16	ALD	V	301	8	-	0/39/39/39	0/1/1/1
16	ALD	Y	301	11	-	0/39/39/39	0/1/1/1
16	ALD	b	201	14	-	0/39/39/39	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
16	b	201	ALD	C7-C3	-4.27	1.40	1.50
16	V	301	ALD	C7-C3	-4.26	1.40	1.50
16	H	301	ALD	C7-C3	-4.16	1.40	1.50
16	K	301	ALD	C7-C3	-4.13	1.41	1.50
16	N	201	ALD	C7-C3	-4.06	1.41	1.50
16	Y	301	ALD	C7-C3	-4.02	1.41	1.50

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
16	N	201	ALD	O31-C9-N10	-3.14	119.47	124.87
16	b	201	ALD	O31-C9-N10	-3.01	119.68	124.87
16	H	301	ALD	O31-C9-N10	-3.01	119.70	124.87
16	V	301	ALD	O31-C9-N10	-2.97	119.76	124.87

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
16	V	301	ALD	C31-C30-C11	-2.68	107.73	115.47
16	Y	301	ALD	O31-C9-N10	-2.67	120.28	124.87
16	K	301	ALD	O31-C9-N10	-2.32	120.88	124.87
16	Y	301	ALD	O8-C9-O31	-2.31	119.58	124.22
16	K	301	ALD	C30-C11-C12	-2.30	104.93	110.60
16	K	301	ALD	C31-C30-C11	-2.12	109.34	115.47
16	H	301	ALD	C31-C30-C11	-2.08	109.48	115.47
16	b	201	ALD	C30-C11-N10	2.29	116.01	110.55
16	N	201	ALD	C30-C11-N10	2.63	116.84	110.55
16	K	301	ALD	O8-C9-N10	3.78	118.67	110.54
16	V	301	ALD	O8-C9-N10	4.23	119.64	110.54
16	H	301	ALD	O8-C9-N10	4.30	119.77	110.54
16	b	201	ALD	O8-C9-N10	4.33	119.86	110.54
16	Y	301	ALD	O8-C9-N10	4.46	120.12	110.54
16	N	201	ALD	O8-C9-N10	4.50	120.22	110.54

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

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%)	-0.61	3 (1%) 79 80	24, 31, 54, 85	0
1	O	250/250 (100%)	-0.64	2 (0%) 86 86	24, 33, 59, 87	0
2	B	244/258 (94%)	-0.45	3 (1%) 79 80	25, 37, 73, 125	0
2	P	244/258 (94%)	-0.49	5 (2%) 65 67	24, 32, 68, 125	0
3	C	240/254 (94%)	-0.49	2 (0%) 86 86	24, 36, 76, 103	0
3	Q	240/254 (94%)	-0.31	10 (4%) 37 39	27, 39, 99, 131	0
4	D	235/260 (90%)	-0.64	0 100 100	26, 34, 53, 78	0
4	R	235/260 (90%)	-0.57	2 (0%) 84 85	29, 42, 67, 105	0
5	E	231/234 (98%)	-0.61	1 (0%) 92 92	27, 43, 65, 97	0
5	S	231/234 (98%)	-0.62	1 (0%) 92 92	24, 37, 61, 88	0
6	F	243/288 (84%)	-0.50	0 100 100	33, 45, 76, 101	0
6	T	243/288 (84%)	-0.59	0 100 100	26, 32, 58, 82	0
7	G	241/252 (95%)	-0.58	1 (0%) 92 92	25, 32, 58, 107	0
7	U	241/252 (95%)	-0.67	2 (0%) 86 86	24, 29, 49, 85	0
8	H	222/232 (95%)	-0.67	1 (0%) 90 91	26, 31, 51, 165	0
8	V	222/232 (95%)	-0.67	0 100 100	27, 34, 52, 117	0
9	I	204/205 (99%)	-0.71	1 (0%) 90 91	25, 32, 50, 66	0
9	W	204/205 (99%)	-0.75	1 (0%) 90 91	25, 29, 45, 62	0
10	J	195/198 (98%)	-0.74	1 (0%) 90 91	24, 29, 44, 89	0
10	X	195/198 (98%)	-0.83	2 (1%) 82 83	24, 27, 39, 79	0
11	K	212/212 (100%)	-0.76	0 100 100	24, 28, 41, 94	0
11	Y	212/212 (100%)	-0.71	0 100 100	25, 33, 55, 84	0
12	L	222/222 (100%)	-0.71	0 100 100	26, 32, 46, 57	0
12	Z	222/222 (100%)	-0.72	0 100 100	26, 34, 53, 66	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
13	M	233/246 (94%)	-0.70	0 100 100	25, 33, 48, 57	0
13	a	233/246 (94%)	-0.78	1 (0%) 92 92	24, 29, 39, 44	0
14	N	196/196 (100%)	-0.71	1 (0%) 90 91	25, 28, 44, 173	0
14	b	196/196 (100%)	-0.77	0 100 100	23, 26, 44, 184	0
All	All	6336/6614 (95%)	-0.64	40 (0%) 89 89	23, 33, 60, 184	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	219	ALA	6.4
2	P	219	ALA	4.9
2	P	51	VAL	4.7
2	P	218	GLY	4.4
3	Q	239	GLN	4.1
3	Q	204	GLY	3.8
1	A	1	MET	3.7
1	O	1	MET	3.7
3	C	206	LYS	3.6
3	Q	238	LYS	3.6
9	I	1	SER	3.6
14	N	1	THR	3.4
3	Q	234	ILE	3.4
3	Q	236	GLN	3.3
3	Q	206	LYS	3.3
5	S	202	ASP	2.8
1	A	250	LEU	2.8
2	B	218	GLY	2.8
2	B	220	ASN	2.8
3	Q	203	THR	2.8
3	C	238	LYS	2.7
8	H	1	THR	2.7
10	J	1	MET	2.7
1	O	249	ALA	2.5
3	Q	51	LYS	2.5
2	P	52	THR	2.5
10	X	194	ASP	2.5
2	P	222	GLY	2.5
7	U	2	GLY	2.5
10	X	1	MET	2.4
3	Q	48	SER	2.4
7	G	242	GLN	2.3

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Mol	Chain	Res	Type	RSRZ
9	W	1	SER	2.2
3	Q	50	LEU	2.2
5	E	202	ASP	2.2
1	A	2	THR	2.1
4	R	125	LEU	2.1
7	U	242	GLN	2.1
4	R	241	ALA	2.0
13	a	1	THR	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
16	ALD	Y	301	34/34	0.80	0.29	6.42	102,131,173,264	0
16	ALD	K	301	34/34	0.80	0.29	5.34	94,114,141,173	0
16	ALD	V	301	34/34	0.87	0.29	4.59	92,128,171,201	0
16	ALD	b	201	34/34	0.88	0.26	3.61	65,109,187,308	0
16	ALD	H	301	34/34	0.86	0.28	3.19	81,137,193,223	0
15	MG	V	302	1/1	0.91	0.16	3.18	110,110,110,110	0
15	MG	K	302	1/1	0.95	0.14	3.03	67,67,67,67	0
16	ALD	N	201	34/34	0.83	0.28	2.83	57,121,178,330	0
15	MG	N	202	1/1	0.92	0.12	0.55	60,60,60,60	0
15	MG	Y	302	1/1	0.94	0.11	0.29	42,42,42,42	0
15	MG	I	301	1/1	0.92	0.10	-0.82	98,98,98,98	0
15	MG	G	301	1/1	0.99	0.07	-1.74	47,47,47,47	0
15	MG	Z	301	1/1	0.97	0.06	-2.61	97,97,97,97	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
15	MG	H	302	1/1	0.97	0.08	-	109,109,109,109	0

6.5 Other polymers [i](#)

There are no such residues in this entry.