



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 08:32 AM UTC

PDB ID : 5LF0 / pdb_00005lf0
Title : Human 20S proteasome complex with Epoxomicin at 2.4 Angstrom
Authors : Schrader, J.; Henneberg, F.; Mata, R.; Tittmann, K.; Schneider, T.R.; Stark, H.; Bourenkov, G.; Chari, A.
Deposited on : 2016-06-30
Resolution : 2.41 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

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

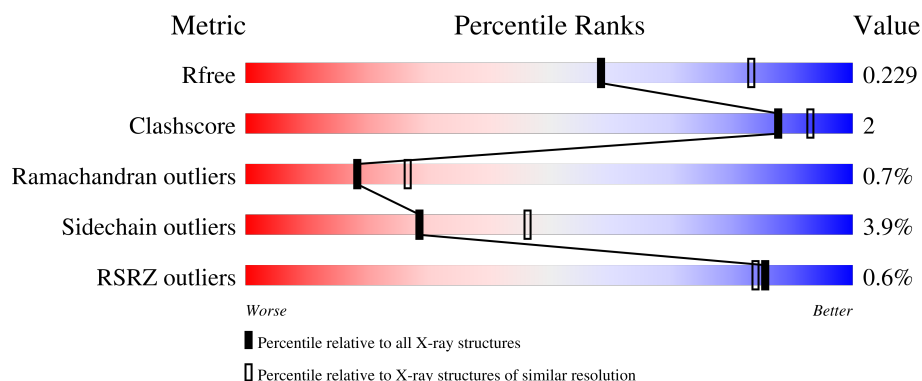
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.41 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}	180053	6062 (2.44-2.40)
Clashscore	190562	6562 (2.44-2.40)
Ramachandran outliers	187476	6481 (2.44-2.40)
Sidechain outliers	187428	6482 (2.44-2.40)
RSRZ outliers	180081	6066 (2.44-2.40)

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

Mol	Chain	Length	Quality of chain
1	A	234	
1	O	234	
2	B	261	
2	P	261	
3	C	248	

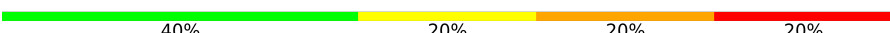
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Mol	Chain	Length	Quality of chain
3	Q	248	
4	D	241	
4	R	241	
5	E	263	
5	S	263	
6	F	255	
6	T	255	
7	G	246	
7	U	246	
8	H	234	
8	V	234	
9	I	205	
9	W	205	
10	J	201	
10	X	201	
11	K	204	
11	Y	204	
12	L	213	
12	Z	213	
13	M	219	
13	a	219	
14	N	205	
14	b	205	
15	c	5	
15	d	5	

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Mol	Chain	Length	Quality of chain
15	e	5	 40%60%
15	f	5	 60%40%
15	g	5	 60%40%
15	h	5	 40%20%20%20%

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	IML	c	2	X	-	-	-
15	IML	d	2	X	-	-	-
15	IML	e	2	X	-	-	-
15	IML	f	2	X	-	-	-
15	IML	g	2	X	-	-	-
15	IML	h	2	X	-	-	-
7	6V1	U	47	X	-	-	-

2 Entry composition

There are 20 unique types of molecules in this entry. The entry contains 51776 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	230	Total	C	N	O	S	0	3	0
			1788	1145	301	336	6			
1	O	230	Total	C	N	O	S	0	0	0
			1741	1111	293	331	6			

- Molecule 2 is a protein called Proteasome subunit alpha type-4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	248	Total	C	N	O	S	0	2	0
			1926	1220	332	363	11			
2	P	248	Total	C	N	O	S	0	2	0
			1909	1206	325	367	11			

- Molecule 3 is a protein called Proteasome subunit alpha type-7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	237	Total	C	N	O	S	0	2	0
			1798	1121	320	352	5			
3	Q	239	Total	C	N	O	S	0	0	0
			1820	1136	320	359	5			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	233	Total	C	N	O	S	0	1	0
			1762	1105	290	356	11			
4	R	233	Total	C	N	O	S	0	1	0
			1753	1103	293	346	11			

- Molecule 5 is a protein called Proteasome subunit alpha type-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	234	Total	C	N	O	S	0	1	0
			1822	1144	325	342	11			
5	S	238	Total	C	N	O	S	0	3	0
			1875	1175	340	349	11			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	148	6V1	CYS	modified residue	UNP P25786
S	148	6V1	CYS	modified residue	UNP P25786

- Molecule 6 is a protein called Proteasome subunit alpha type-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	F	239	Total	C	N	O	S	0	4	0
			1888	1198	325	353	12			
6	T	240	Total	C	N	O	S	0	1	0
			1856	1178	315	351	12			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	G	244	Total	C	N	O	S	0	2	0
			1912	1214	321	364	13			
7	U	238	Total	C	N	O	S	0	1	0
			1815	1147	304	350	14			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	47	6V1	CYS	modified residue	UNP P60900
G	161	6V1	CYS	modified residue	UNP P60900
U	47	6V1	CYS	modified residue	UNP P60900
U	161	6V1	CYS	modified residue	UNP P60900

- Molecule 8 is a protein called Proteasome subunit beta type-7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	H	220	Total	C	N	O	S	0	2	0
			1664	1047	284	320	13			
8	V	220	Total	C	N	O	S	0	2	0
			1622	1023	269	318	12			

- 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	3	0
			1613	1028	270	295	20			
9	W	204	Total	C	N	O	S	0	2	0
			1599	1018	267	295	19			

- Molecule 10 is a protein called Proteasome subunit beta type-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	J	196	Total	C	N	O	S	0	3	0
			1590	1021	271	288	10			
10	X	196	Total	C	N	O	S	0	2	0
			1572	1012	266	284	10			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
J	91	6V1	CYS	modified residue	UNP P49721
X	91	6V1	CYS	modified residue	UNP P49721

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
11	K	200	Total	C	N	O	S	0	0	0
			1545	974	269	293	9			
11	Y	201	Total	C	N	O	S	0	3	0
			1580	996	280	294	10			

- Molecule 12 is a protein called Proteasome subunit beta type-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
12	L	213	Total	C	N	O	S	0	2	0
			1636	1038	277	310	11			
12	Z	213	Total	C	N	O	S	0	1	0
			1642	1041	280	310	11			

- Molecule 13 is a protein called Proteasome subunit beta type-4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
13	M	216	Total	C	N	O	S	0	1	0
			1692	1067	291	322	12			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
13	a	216	Total	C	N	O	S	0	2	0
			1688	1064	291	321	12			

- Molecule 14 is a protein called Proteasome subunit beta type-6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
14	N	202	Total	C	N	O	S	0	1	0
			1519	953	258	295	13			
14	b	203	Total	C	N	O	S	0	1	0
			1524	956	259	296	13			

- Molecule 15 is a protein called EPOXOMICIN (peptide inhibitor).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
15	c	5	Total	C	N	O	0	0	0
			39	28	4	7			
15	d	5	Total	C	N	O	0	0	0
			39	28	4	7			
15	e	5	Total	C	N	O	0	0	0
			39	28	4	7			
15	f	5	Total	C	N	O	0	0	0
			39	28	4	7			
15	g	5	Total	C	N	O	0	0	0
			39	28	4	7			
15	h	5	Total	C	N	O	0	0	0
			39	28	4	7			

- Molecule 16 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
16	A	4	Total	Cl	0	0
			4	4		
16	B	2	Total	Cl	0	0
			2	2		
16	C	2	Total	Cl	0	0
			2	2		
16	D	1	Total	Cl	0	0
			1	1		
16	E	4	Total	Cl	0	0
			4	4		
16	F	1	Total	Cl	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
16	G	2	Total 2	Cl 2	0	0
16	H	1	Total 1	Cl 1	0	0
16	I	1	Total 1	Cl 1	0	0
16	K	4	Total 4	Cl 4	0	0
16	M	3	Total 3	Cl 3	0	0
16	N	4	Total 4	Cl 4	0	0
16	O	4	Total 4	Cl 4	0	0
16	P	1	Total 1	Cl 1	0	0
16	Q	2	Total 2	Cl 2	0	0
16	R	2	Total 2	Cl 2	0	0
16	S	3	Total 3	Cl 3	0	0
16	U	1	Total 1	Cl 1	0	0
16	V	1	Total 1	Cl 1	0	0
16	W	1	Total 1	Cl 1	0	0
16	Y	5	Total 5	Cl 5	0	0
16	a	3	Total 3	Cl 3	0	0
16	b	4	Total 4	Cl 4	0	0
16	c	1	Total 1	Cl 1	0	0
16	f	1	Total 1	Cl 1	0	0

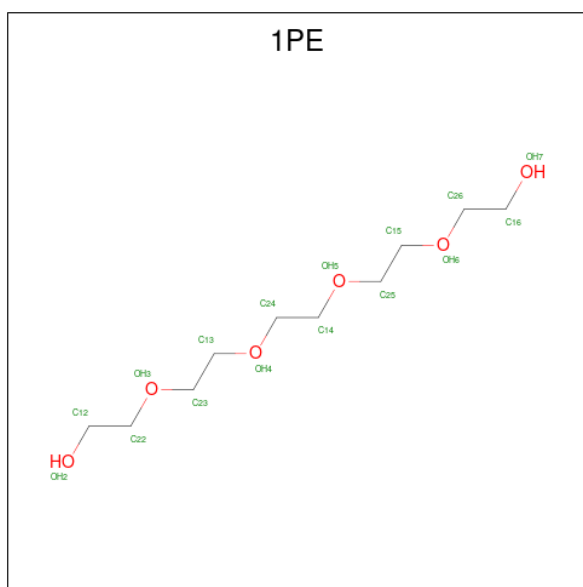
- Molecule 17 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
17	G	1	Total K 1 1	0	0
17	L	1	Total K 1 1	0	0
17	N	1	Total K 1 1	0	0
17	U	1	Total K 1 1	0	0
17	Z	1	Total K 1 1	0	0
17	b	1	Total K 1 1	0	0

- Molecule 18 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
18	H	2	Total Mg 2 2	0	0
18	I	1	Total Mg 1 1	0	0
18	J	1	Total Mg 1 1	0	0
18	K	1	Total Mg 1 1	0	0
18	V	2	Total Mg 2 2	0	0
18	W	1	Total Mg 1 1	0	0
18	X	1	Total Mg 1 1	0	0
18	Y	1	Total Mg 1 1	0	0

- Molecule 19 is PENTAETHYLENE GLYCOL (CCD ID: 1PE) (formula: C₁₀H₂₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
19	H	1	Total	C	O	0	0
			16	10	6		
19	I	1	Total	C	O	0	0
			16	10	6		
19	K	1	Total	C	O	0	0
			16	10	6		
19	L	1	Total	C	O	0	0
			16	10	6		
19	N	1	Total	C	O	0	0
			16	10	6		
19	W	1	Total	C	O	0	0
			16	10	6		
19	Z	1	Total	C	O	0	0
			16	10	6		
19	a	1	Total	C	O	0	0
			16	10	6		
19	b	1	Total	C	O	0	0
			16	10	6		

- Molecule 20 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
20	A	100	Total	O	0	0
			100	100		
20	B	116	Total	O	0	0
			116	116		
20	C	62	Total	O	0	0
			62	62		

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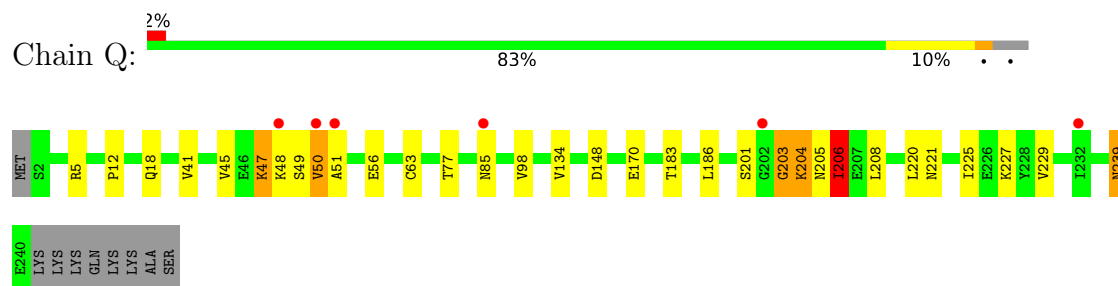
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
20	D	81	Total 81	O 81	0	0
20	E	127	Total 127	O 127	0	0
20	F	164	Total 164	O 164	0	0
20	G	172	Total 172	O 172	0	0
20	H	144	Total 144	O 144	0	0
20	I	142	Total 142	O 142	0	0
20	J	119	Total 119	O 119	0	0
20	K	95	Total 95	O 95	0	0
20	L	114	Total 114	O 114	0	0
20	M	141	Total 141	O 141	0	0
20	N	143	Total 143	O 143	0	0
20	O	71	Total 71	O 71	0	0
20	P	102	Total 102	O 102	0	0
20	Q	57	Total 57	O 57	0	0
20	R	109	Total 109	O 109	0	0
20	S	101	Total 101	O 101	0	0
20	T	73	Total 73	O 73	0	0
20	U	80	Total 80	O 80	0	0
20	V	104	Total 104	O 104	0	0
20	W	89	Total 89	O 89	0	0
20	X	109	Total 109	O 109	0	0

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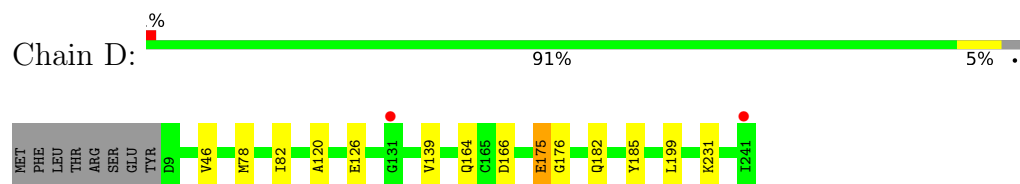
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
20	Y	135	Total 135	O 135	0	0
20	Z	155	Total 155	O 155	0	0
20	a	154	Total 154	O 154	0	0
20	b	110	Total 110	O 110	0	0
20	c	1	Total 1	O 1	0	0
20	d	1	Total 1	O 1	0	0
20	g	1	Total 1	O 1	0	0
20	h	1	Total 1	O 1	0	0

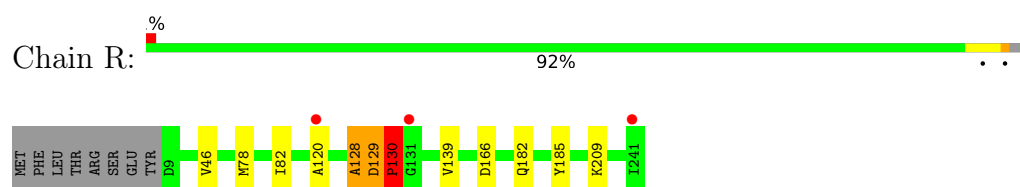
- Molecule 3: Proteasome subunit alpha type-7



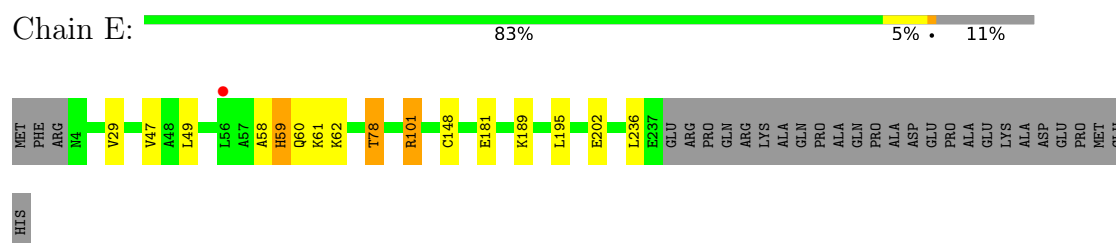
- Molecule 4: Proteasome subunit alpha type-5



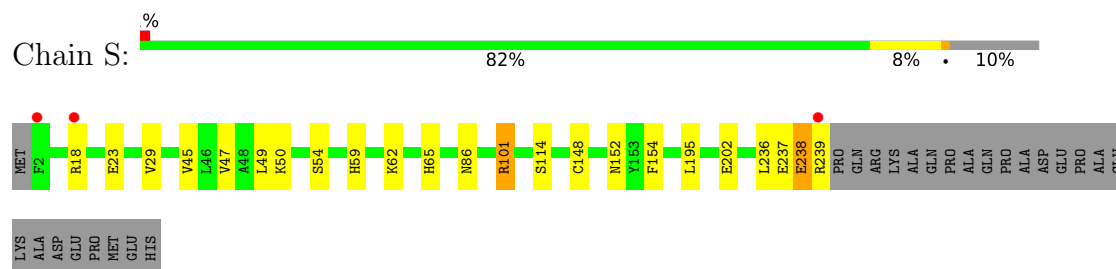
- Molecule 4: Proteasome subunit alpha type-5



- Molecule 5: Proteasome subunit alpha type-1

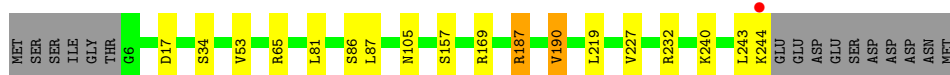


- Molecule 5: Proteasome subunit alpha type-1

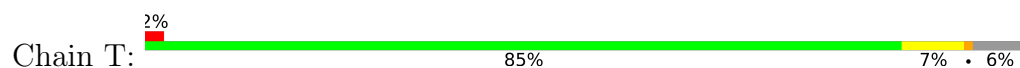


- Molecule 6: Proteasome subunit alpha type-3





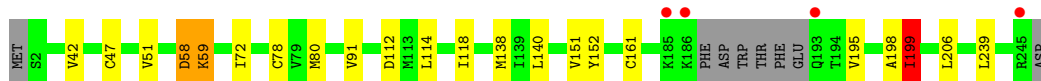
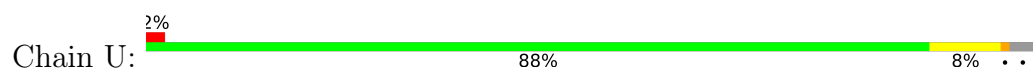
- Molecule 6: Proteasome subunit alpha type-3



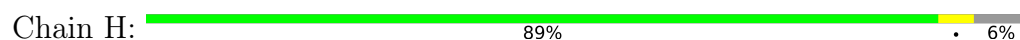
- Molecule 7: Proteasome subunit alpha type-6



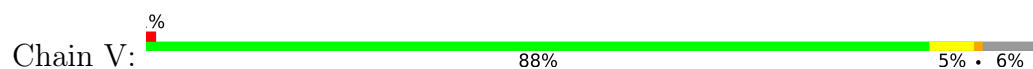
- Molecule 7: Proteasome subunit alpha type-6



- Molecule 8: Proteasome subunit beta type-7



- Molecule 8: Proteasome subunit beta type-7



- Molecule 9: Proteasome subunit beta type-3



- Molecule 9: Proteasome subunit beta type-3





- Molecule 10: Proteasome subunit beta type-2

Chain J: 89% 7% ..



- Molecule 10: Proteasome subunit beta type-2

Chain X: 89% 6% ..



- Molecule 11: Proteasome subunit beta type-5

Chain K: 88% 9% .



- Molecule 11: Proteasome subunit beta type-5

Chain Y: 89% 8% ..



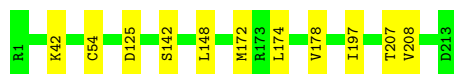
- Molecule 12: Proteasome subunit beta type-1

Chain L: 94% 6%



- Molecule 12: Proteasome subunit beta type-1

Chain Z: 95% 5%



- Molecule 13: Proteasome subunit beta type-4

Chain M: 93% 6% .



- Molecule 13: Proteasome subunit beta type-4

Chain a: 94% 5% ..



- Molecule 14: Proteasome subunit beta type-6

Chain N: 95% ..



- Molecule 14: Proteasome subunit beta type-6

Chain b: 93% 5% ..



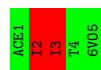
- Molecule 15: EPOXOMICIN (peptide inhibitor)

Chain c: 60% 20% 20%



- Molecule 15: EPOXOMICIN (peptide inhibitor)

Chain d: 60% 40%



- Molecule 15: EPOXOMICIN (peptide inhibitor)

Chain e: 40% 60%



- Molecule 15: EPOXOMICIN (peptide inhibitor)

Chain f: 60% 40%



- Molecule 15: EPOXOMICIN (peptide inhibitor)



- Molecule 15: EPOXOMICIN (peptide inhibitor)



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	113.07Å 202.37Å 314.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	170.23 – 2.41 170.23 – 2.41	Depositor EDS
% Data completeness (in resolution range)	99.5 (170.23-2.41) 99.5 (170.23-2.41)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.24 (at 2.43Å)	Xtriage
Refinement program	REFMAC 5.8.0103	Depositor
R, R_{free}	0.179 , 0.229 0.182 , 0.229	Depositor DCC
R_{free} test set	13832 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	52.7	Xtriage
Anisotropy	0.278	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 48.2	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	0.96	EDS
Total number of atoms	51776	wwPDB-VP
Average B, all atoms (Å ²)	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.91% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: CL, ACE, 1PE, IML, MG, YCM, K, 6VO, 6V1

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	0.72	0/1833	0.90	1/2489 (0.0%)
1	O	0.69	0/1778	0.89	0/2419
2	B	0.72	0/1962	0.90	1/2649 (0.0%)
2	P	0.71	0/1945	0.92	1/2631 (0.0%)
3	C	0.77	0/1818	1.00	1/2469 (0.0%)
3	Q	0.78	1/1834 (0.1%)	1.02	5/2490 (0.2%)
4	D	0.73	0/1789	0.93	1/2424 (0.0%)
4	R	0.76	0/1780	0.98	2/2408 (0.1%)
5	E	0.75	0/1842	0.91	1/2493 (0.0%)
5	S	0.72	0/1901	0.93	0/2571
6	F	0.73	0/1935	0.91	2/2605 (0.1%)
6	T	0.72	0/1894	0.97	6/2556 (0.2%)
7	G	0.74	0/1909	0.92	4/2579 (0.2%)
7	U	0.72	0/1804	0.91	1/2441 (0.0%)
8	H	0.75	1/1697 (0.1%)	0.91	0/2299
8	V	0.70	1/1655 (0.1%)	0.87	0/2251
9	I	0.70	0/1648	1.01	6/2219 (0.3%)
9	W	0.66	1/1630 (0.1%)	0.91	6/2197 (0.3%)
10	J	0.71	0/1613	0.85	0/2180
10	X	0.71	0/1595	0.85	0/2157
11	K	0.73	0/1576	0.91	0/2131
11	Y	0.75	0/1620	0.92	3/2185 (0.1%)
12	L	0.69	0/1672	0.84	0/2257
12	Z	0.74	0/1675	0.86	0/2257
13	M	0.73	0/1728	0.89	0/2339
13	a	0.71	0/1724	0.88	1/2336 (0.0%)
14	N	0.76	0/1548	0.91	0/2095
14	b	0.71	0/1554	0.90	1/2104 (0.0%)
15	c	0.87	0/14	2.00	1/18 (5.6%)
15	d	0.54	0/14	2.06	1/18 (5.6%)
15	e	0.83	0/14	1.66	0/18
15	f	0.89	0/14	2.06	0/18

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
15	g	0.87	0/14	2.33	2/18 (11.1%)
15	h	0.69	0/14	2.41	1/18 (5.6%)
All	All	0.73	4/49043 (0.0%)	0.92	48/66339 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	P	0	4
3	C	0	1
3	Q	0	2
4	D	0	2
4	R	0	1
7	U	1	0
9	I	0	1
9	W	0	1
10	J	0	2
10	X	0	1
11	Y	0	1
13	a	0	1
15	c	1	0
15	d	2	0
15	e	1	0
15	f	1	0
15	g	2	0
15	h	2	0
All	All	10	17

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	W	77	GLU	CD-OE2	5.92	1.36	1.25
3	Q	206	ILE	CA-C	5.72	1.60	1.52
8	V	220	GLU	C-O	5.47	1.34	1.23
8	H	47	GLY	C-O	-5.08	1.21	1.24

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	I	16[A]	LYS	CA-C-N	10.24	136.72	122.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	I	16[A]	LYS	C-N-CA	10.24	136.72	122.07
9	I	16[B]	LYS	CA-C-N	10.24	136.72	122.07
9	I	16[B]	LYS	C-N-CA	10.24	136.72	122.07
6	T	6	GLY	N-CA-C	9.40	135.47	113.18
2	P	246	LYS	N-CA-C	7.80	120.87	111.82
4	R	120[A]	ALA	N-CA-C	-7.61	99.02	110.28
4	R	120[B]	ALA	N-CA-C	-7.61	99.02	110.28
6	T	190	VAL	CB-CA-C	-7.54	102.17	112.04
6	F	190	VAL	CB-CA-C	-7.38	102.38	112.04
9	I	79	ARG	N-CA-C	7.31	119.27	108.60
6	T	6	GLY	CA-C-N	6.99	134.28	121.70
6	T	6	GLY	C-N-CA	6.99	134.28	121.70
9	W	79	ARG	N-CA-C	6.94	118.73	108.60
4	D	120	ALA	N-CA-C	6.85	122.67	111.37
15	d	3	ILE	CA-CB-CG2	6.74	121.96	110.50
3	Q	12	PRO	N-CA-C	6.43	122.20	113.84
3	Q	220	LEU	CA-C-N	6.33	133.10	121.70
3	Q	220	LEU	C-N-CA	6.33	133.10	121.70
7	G	244	GLU	N-CA-C	6.03	118.25	110.65
5	E	236	LEU	N-CA-C	-5.93	103.18	110.65
1	A	221	THR	CB-CA-C	5.70	117.07	108.86
3	Q	220	LEU	N-CA-C	-5.64	100.52	109.76
15	g	3	ILE	CA-CB-CG2	5.52	119.88	110.50
7	G	183	VAL	CB-CA-C	-5.48	103.66	112.16
9	W	16[A]	LYS	CA-C-N	5.45	133.43	123.13
9	W	16[A]	LYS	C-N-CA	5.45	133.43	123.13
9	W	16[B]	LYS	CA-C-N	5.45	133.43	123.13
9	W	16[B]	LYS	C-N-CA	5.45	133.43	123.13
11	Y	1	THR	N-CA-CB	5.43	120.73	111.50
15	h	3	ILE	CB-CA-C	5.38	122.36	111.60
7	G	183	VAL	N-CA-C	5.27	117.64	111.05
14	b	22	THR	CB-CA-C	-5.27	102.25	110.74
6	T	190	VAL	N-CA-CB	5.26	117.05	110.47
3	Q	204	LYS	N-CA-C	5.24	115.97	107.32
15	c	3	ILE	CA-CB-CG2	5.24	119.40	110.50
15	g	3	ILE	CB-CG1-CD1	-5.23	102.82	113.80
13	a	53	ALA	N-CA-C	5.22	116.90	108.34
9	I	193	LYS	N-CA-C	5.21	116.44	108.42
6	F	190	VAL	N-CA-CB	5.19	116.96	110.47
3	C	47	LYS	N-CA-C	5.14	121.74	110.80
2	B	55	LEU	N-CA-C	5.11	119.38	113.20
7	U	199	ILE	CB-CA-C	5.10	120.48	112.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	T	143	ASN	N-CA-C	5.10	116.92	111.36
7	G	183	VAL	N-CA-CB	5.09	119.36	110.65
9	W	193	LYS	N-CA-C	5.07	116.23	108.42
11	Y	199	TYR	CA-C-N	5.04	130.78	121.70
11	Y	199	TYR	C-N-CA	5.04	130.78	121.70

All (10) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
7	U	47	6V1	C1
15	c	2	IML	CB
15	d	2	IML	CB,CA
15	e	2	IML	CA
15	f	2	IML	CB
15	g	2	IML	CB,CA
15	h	2	IML	CB,CA

All (17) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	C	237	GLU	Peptide
4	D	175[A]	GLU	Peptide
4	D	175[B]	GLU	Peptide
9	I	78	GLY	Peptide
10	J	1[A]	MET	Peptide
10	J	1[B]	MET	Peptide
2	P	244	GLU	Peptide
2	P	245	ALA	Peptide
2	P	53	HIS	Peptide
2	P	54	LYS	Peptide
3	Q	47	LYS	Peptide
3	Q	49	SER	Peptide
4	R	130	PRO	Peptide
9	W	78	GLY	Peptide
10	X	1	MET	Peptide
11	Y	199	TYR	Peptide
13	a	215	ILE	Peptide

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	1788	0	1761	6	0
1	O	1741	0	1683	4	0
2	B	1926	0	1924	3	0
2	P	1909	0	1874	10	0
3	C	1798	0	1718	11	0
3	Q	1820	0	1749	12	0
4	D	1762	0	1709	3	0
4	R	1753	0	1726	5	0
5	E	1822	0	1779	7	0
5	S	1875	0	1818	14	0
6	F	1888	0	1882	7	0
6	T	1856	0	1816	6	0
7	G	1912	0	1882	4	0
7	U	1815	0	1748	11	0
8	H	1664	0	1678	4	0
8	V	1622	0	1592	3	0
9	I	1613	0	1646	7	0
9	W	1599	0	1621	7	0
10	J	1590	0	1581	10	0
10	X	1572	0	1559	11	0
11	K	1545	0	1495	11	0
11	Y	1580	0	1555	13	0
12	L	1636	0	1625	6	0
12	Z	1642	0	1635	4	0
13	M	1692	0	1670	6	0
13	a	1688	0	1658	3	0
14	N	1519	0	1493	6	0
14	b	1524	0	1493	7	0
15	c	39	0	34	3	0
15	d	39	0	34	4	0
15	e	39	0	34	0	0
15	f	39	0	34	2	0
15	g	39	0	34	4	0
15	h	39	0	34	2	0
16	A	4	0	0	0	0
16	B	2	0	0	1	0
16	C	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
16	D	1	0	0	0	0
16	E	4	0	0	0	0
16	F	1	0	0	0	0
16	G	2	0	0	1	0
16	H	1	0	0	0	0
16	I	1	0	0	0	0
16	K	4	0	0	1	0
16	M	3	0	0	0	0
16	N	4	0	0	0	0
16	O	4	0	0	0	0
16	P	1	0	0	0	0
16	Q	2	0	0	0	0
16	R	2	0	0	0	0
16	S	3	0	0	0	0
16	U	1	0	0	0	0
16	V	1	0	0	0	0
16	W	1	0	0	0	0
16	Y	5	0	0	0	0
16	a	3	0	0	1	0
16	b	4	0	0	1	0
16	c	1	0	0	0	0
16	f	1	0	0	0	0
17	G	1	0	0	0	0
17	L	1	0	0	0	0
17	N	1	0	0	0	0
17	U	1	0	0	0	0
17	Z	1	0	0	0	0
17	b	1	0	0	0	0
18	H	2	0	0	0	0
18	I	1	0	0	0	0
18	J	1	0	0	0	0
18	K	1	0	0	0	0
18	V	2	0	0	0	0
18	W	1	0	0	0	0
18	X	1	0	0	0	0
18	Y	1	0	0	0	0
19	H	16	0	22	0	0
19	I	16	0	22	0	0
19	K	16	0	22	0	0
19	L	16	0	22	0	0
19	N	16	0	22	0	0
19	W	16	0	22	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
19	Z	16	0	22	0	0
19	a	16	0	22	0	0
19	b	16	0	22	0	0
20	A	100	0	0	1	0
20	B	116	0	0	0	0
20	C	62	0	0	0	0
20	D	81	0	0	0	0
20	E	127	0	0	1	0
20	F	164	0	0	3	0
20	G	172	0	0	0	0
20	H	144	0	0	0	0
20	I	142	0	0	0	0
20	J	119	0	0	1	0
20	K	95	0	0	0	0
20	L	114	0	0	1	0
20	M	141	0	0	0	0
20	N	143	0	0	0	0
20	O	71	0	0	1	0
20	P	102	0	0	0	0
20	Q	57	0	0	0	0
20	R	109	0	0	1	0
20	S	101	0	0	4	0
20	T	73	0	0	0	0
20	U	80	0	0	0	0
20	V	104	0	0	1	0
20	W	89	0	0	0	0
20	X	109	0	0	0	0
20	Y	135	0	0	0	0
20	Z	155	0	0	1	0
20	a	154	0	0	0	0
20	b	110	0	0	0	0
20	c	1	0	0	0	0
20	d	1	0	0	0	0
20	g	1	0	0	0	0
20	h	1	0	0	0	0
All	All	51776	0	47772	189	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 (189) 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)
12:L:144:MET:HE1	12:L:185:ARG:HB2	1.57	0.87
8:H:48:THR:HG23	15:c:2:IML:HD11	1.54	0.86
6:F:105:ASN:ND2	20:F:401:HOH:O	2.16	0.77
11:K:35:ILE:HD11	11:K:45:MET:SD	2.33	0.69
2:P:12:PHE:H	3:Q:18:GLN:HE22	1.42	0.68
9:I:25[A]:ARG:HG2	9:I:25[A]:ARG:HH11	1.59	0.67
5:E:47:VAL:HG12	5:E:195:LEU:HD22	1.76	0.67
11:Y:35:ILE:HD11	11:Y:45:MET:SD	2.34	0.67
3:Q:204:LYS:HA	3:Q:205:ASN:O	1.95	0.67
11:Y:199:TYR:HA	11:Y:200:SER:HB2	1.76	0.66
1:A:108:GLN:HE21	1:A:112:ARG:HH12	1.42	0.66
5:S:47:VAL:HG12	5:S:195:LEU:HD22	1.76	0.66
9:W:25[A]:ARG:HH11	9:W:25[A]:ARG:HG2	1.61	0.65
11:K:141:ARG:NH1	10:X:166:GLU:OE2	2.29	0.65
15:d:2:IML:HG22	15:d:2:IML:HN3	1.77	0.65
6:F:227:VAL:O	6:F:232[B]:ARG:NH1	2.28	0.64
15:g:2:IML:HG22	15:g:2:IML:HN3	1.80	0.63
1:A:88[B]:ARG:NH2	20:A:401:HOH:O	2.29	0.62
5:S:101:ARG:NH1	20:S:401:HOH:O	2.32	0.61
9:I:124:ASP:HB3	15:c:3:ILE:HG23	1.82	0.61
14:N:35:THR:CG2	14:N:45:ARG:HE	2.14	0.60
15:f:2:IML:HN2	15:f:2:IML:HG12	1.83	0.60
5:E:58:ALA:O	5:E:59:HIS:CB	2.50	0.59
10:J:1[A]:MET:HE1	10:J:134:TYR:H	1.66	0.59
10:J:166:GLU:OE2	11:Y:141[A]:ARG:NH1	2.35	0.58
7:U:118:ILE:HG13	7:U:138:MET:HE1	1.85	0.58
12:L:144:MET:CE	12:L:185:ARG:HB2	2.31	0.57
10:X:1:MET:HE1	10:X:134:TYR:H	1.70	0.57
1:A:51:GLN:HE22	1:A:58[B]:GLU:HB3	1.69	0.56
11:Y:20:ALA:HB1	15:g:3:ILE:CD1	2.34	0.56
10:J:101:ASN:HD22	10:J:119:ASP:HA	1.71	0.56
5:S:18[A]:ARG:HD2	5:S:23:GLU:OE2	2.06	0.56
8:V:213:THR:HB	9:W:198:THR:OG1	2.06	0.56
11:Y:199:TYR:HA	11:Y:200:SER:CB	2.36	0.56
16:B:301:CL:CL	16:B:302:CL:CL	2.98	0.55
14:b:35:THR:CG2	14:b:45:ARG:HE	2.18	0.55
7:U:58:ASP:O	7:U:59:LYS:CB	2.55	0.55
7:U:80[A]:MET:HE2	7:U:91:VAL:HG23	1.90	0.54
3:C:35:VAL:HG13	3:C:191:VAL:CG2	2.36	0.54
10:J:1[A]:MET:HE1	10:J:134:TYR:N	2.22	0.54
4:R:129:ASP:CB	4:R:130:PRO:CD	2.86	0.54
2:P:25[B]:MET:HE3	2:P:25[B]:MET:HA	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:L:125:ASP:HB2	15:d:2:IML:HG13	1.89	0.53
7:G:186:LYS:NZ	16:G:302:CL:CL	2.78	0.53
5:S:49:LEU:O	5:S:62:LYS:HD2	2.09	0.53
5:E:101[A]:ARG:NH1	20:E:401:HOH:O	2.42	0.52
11:Y:52:CYS:SG	11:Y:97[A]:MET:HG3	2.49	0.52
10:X:1:MET:HE1	10:X:134:TYR:N	2.24	0.52
5:E:49:LEU:O	5:E:62:LYS:HD2	2.09	0.52
4:R:78:MET:HG3	4:R:82:ILE:HD12	1.92	0.52
6:F:219:LEU:HD22	20:F:426:HOH:O	2.10	0.51
14:b:14:LEU:HD23	14:b:44:CYS:SG	2.49	0.51
11:K:52:CYS:SG	11:K:97:MET:HG3	2.50	0.51
5:S:238:GLU:CB	5:S:239:ARG:C	2.84	0.51
5:S:50:LYS:HB3	5:S:59:HIS:HB3	1.91	0.51
13:a:5:MET:HE2	14:b:116:MET:HB3	1.93	0.51
5:S:237:GLU:O	5:S:238:GLU:CB	2.58	0.51
3:C:85[B]:ASN:OD1	10:J:70:ARG:NH2	2.43	0.51
14:N:14:LEU:HD23	14:N:44:CYS:SG	2.51	0.50
7:U:195:VAL:O	7:U:199:ILE:HG23	2.11	0.50
6:T:87:LEU:HD13	6:T:131:PHE:CE1	2.46	0.50
15:g:2:IML:HG22	15:g:2:IML:CN	2.41	0.50
11:Y:40:TYR:CD2	11:Y:73:ARG:CZ	2.95	0.50
7:G:72:ILE:HG21	7:G:114:LEU:HD21	1.94	0.49
9:I:13[B]:MET:HE2	9:I:153:TRP:HB2	1.94	0.49
6:T:205:LYS:O	6:T:206:ASP:CG	2.54	0.49
10:X:95:ARG:HB2	10:X:95:ARG:HH11	1.77	0.49
13:M:208:ASN:O	14:b:29:ARG:NH2	2.45	0.49
13:M:5:MET:HE2	14:N:116:MET:HB3	1.94	0.49
7:U:72:ILE:HG21	7:U:114:LEU:HD21	1.93	0.49
10:X:1:MET:HE3	10:X:1:MET:C	2.37	0.49
11:K:40:TYR:CD2	11:K:73:ARG:CZ	2.95	0.48
13:M:27:LEU:HD22	13:M:184:TYR:HB2	1.95	0.48
8:H:132:LEU:HD22	14:N:25:TYR:CE2	2.47	0.48
20:L:501:HOH:O	8:V:195:LYS:CB	2.60	0.48
6:T:24:GLU:HA	6:T:27:MET:HE3	1.96	0.48
12:L:148:LEU:HD23	12:L:178:VAL:CG1	2.43	0.48
12:Z:125:ASP:HB2	15:g:2:IML:HG13	1.94	0.48
15:d:2:IML:HN3	15:d:2:IML:CG2	2.42	0.48
9:W:13:MET:HE1	9:W:165:THR:HB	1.96	0.48
4:D:78:MET:HG3	4:D:82:ILE:HD12	1.96	0.48
2:B:14:PRO:HA	3:C:21:TYR:CD1	2.49	0.48
2:P:246:LYS:O	2:P:249:ARG:NH2	2.46	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:Z:148:LEU:HD23	12:Z:178:VAL:CG1	2.43	0.48
10:J:88:LEU:HB3	10:J:122:ALA:HB2	1.96	0.47
10:X:88:LEU:HB3	10:X:122:ALA:HB2	1.94	0.47
6:F:169[A]:ARG:NH1	20:F:403:HOH:O	2.47	0.47
5:S:86:ASN:ND2	20:S:402:HOH:O	2.48	0.47
11:Y:35:ILE:CD1	11:Y:45:MET:SD	3.02	0.47
13:a:27:LEU:HD22	13:a:184:TYR:HB2	1.96	0.47
1:A:58[B]:GLU:H	1:A:58[B]:GLU:CD	2.22	0.47
9:I:13[A]:MET:HE1	9:I:165:THR:HB	1.96	0.47
7:U:199:ILE:HD11	7:U:239:LEU:HD23	1.96	0.47
11:K:20:ALA:HB2	11:K:31:VAL:HG21	1.97	0.47
3:Q:85:ASN:OD1	10:X:70:ARG:CZ	2.62	0.47
11:K:35:ILE:CD1	11:K:45:MET:SD	3.01	0.46
2:P:246:LYS:N	2:P:246:LYS:HE3	2.31	0.46
14:N:35:THR:HG23	14:N:45:ARG:HE	1.79	0.46
12:Z:172:MET:HE1	12:Z:197:ILE:HD11	1.96	0.46
7:U:51:VAL:HG12	7:U:198:ALA:HB1	1.98	0.46
2:P:155:ASN:OD1	3:Q:77:THR:OG1	2.28	0.46
11:Y:12:VAL:HG13	11:Y:179:VAL:HB	1.97	0.46
2:P:8:ARG:NH2	3:Q:5:ARG:HD2	2.31	0.46
11:Y:20:ALA:HB2	11:Y:31:VAL:HG21	1.97	0.45
11:K:12:VAL:HG13	11:K:179:VAL:HB	1.97	0.45
12:L:172:MET:HE1	12:L:197:ILE:HD11	1.99	0.45
3:Q:85:ASN:OD1	10:X:70:ARG:NH2	2.50	0.45
3:Q:183:THR:CG2	3:Q:186:LEU:HD13	2.47	0.45
1:O:110:VAL:HG22	1:O:135:ILE:HD12	1.98	0.45
5:S:152[B]:ASN:ND2	20:S:403:HOH:O	2.49	0.45
1:A:110:VAL:HG22	1:A:135:ILE:HD12	1.99	0.45
11:K:83:LEU:HD21	11:K:99:THR:HG21	1.99	0.45
3:C:33:VAL:HG21	3:C:195:LEU:HD12	1.98	0.44
20:V:413:HOH:O	15:h:3:ILE:HD11	2.17	0.44
11:K:20:ALA:HB1	15:d:3:ILE:HD13	1.99	0.44
3:Q:203:GLY:HA2	3:Q:229:VAL:HG11	1.99	0.44
13:a:96:MET:HE3	13:a:127:MET:HA	1.99	0.44
6:F:34:SER:OG	6:F:65:ARG:NH1	2.49	0.44
7:U:78:CYS:HB3	7:U:140:LEU:HD23	1.98	0.44
2:P:44:LEU:HD22	2:P:190:LEU:HD13	2.00	0.44
2:B:44:LEU:HD22	2:B:190:LEU:HD13	2.00	0.44
6:F:157:SER:O	7:G:88:ARG:NH2	2.50	0.44
2:P:232:GLU:O	2:P:236:LEU:HD13	2.18	0.44
14:b:35:THR:HG23	14:b:45:ARG:HE	1.81	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:X:46[B]:CYS:SG	10:X:102:LEU:HD22	2.58	0.44
5:E:47:VAL:CG1	5:E:195:LEU:HD22	2.48	0.44
4:D:166:ASP:HB3	4:D:185:TYR:CZ	2.53	0.43
7:G:78:CYS:HB3	7:G:140:LEU:HD23	1.99	0.43
4:R:166:ASP:HB3	4:R:185:TYR:CZ	2.53	0.43
9:W:124:ASP:OD1	9:W:124:ASP:C	2.61	0.43
3:C:47:LYS:CB	3:C:48:LYS:C	2.91	0.43
11:Y:83:LEU:HD21	11:Y:99:THR:HG21	1.99	0.43
15:h:3:ILE:HD13	15:h:3:ILE:HG21	1.83	0.43
11:K:99:THR:HG22	11:K:114:VAL:O	2.19	0.43
4:D:164:GLN:OE1	5:E:58:ALA:HB2	2.18	0.43
8:H:20:ALA:HB1	15:c:3:ILE:CD1	2.49	0.43
13:M:96:MET:HE3	13:M:127:MET:HA	2.00	0.42
2:P:201:MET:HE1	2:P:211:VAL:HG12	2.00	0.42
8:V:132:LEU:HD22	14:b:25:TYR:CZ	2.53	0.42
3:C:203:GLY:CA	3:C:204:LYS:CB	2.97	0.42
5:S:50:LYS:NZ	20:S:404:HOH:O	2.51	0.42
12:Z:42:LYS:HB3	12:Z:54:CYS:O	2.19	0.42
2:B:45:LEU:HD13	2:B:75:SER:HB2	2.01	0.42
1:O:108:GLN:HG3	9:W:77:GLU:OE1	2.19	0.42
10:X:1:MET:HE3	10:X:2:GLU:N	2.34	0.42
6:T:34:SER:OG	6:T:65:ARG:NH1	2.48	0.42
5:E:60:GLN:NE2	5:E:78:THR:HG21	2.34	0.42
5:S:154:PHE:HD2	6:T:63:ASN:HD21	1.68	0.42
1:A:108:GLN:NE2	1:A:112:ARG:HH12	2.15	0.42
6:F:187:ARG:HD2	6:F:187:ARG:HA	1.83	0.42
11:K:32:LYS:NZ	16:K:303:CL:CL	2.90	0.42
1:O:10:THR:HB	1:O:20:GLN:HB2	2.01	0.42
7:U:80[A]:MET:HE2	7:U:91:VAL:CG2	2.48	0.42
3:C:85[B]:ASN:OD1	10:J:70:ARG:CZ	2.67	0.42
7:U:138:MET:HE3	7:U:140:LEU:HD11	2.01	0.42
5:S:47:VAL:CG1	5:S:195:LEU:HD22	2.47	0.42
9:W:13:MET:HE1	9:W:166:ILE:N	2.34	0.42
3:Q:50:VAL:O	3:Q:51:ALA:HB3	2.20	0.42
4:R:129:ASP:CB	4:R:130:PRO:HD3	2.49	0.42
8:H:132:LEU:HD22	14:N:25:TYR:CZ	2.55	0.41
5:S:50:LYS:CB	5:S:59:HIS:HB3	2.49	0.41
11:Y:66:TYR:CD2	11:Y:66:TYR:C	2.98	0.41
12:L:42:LYS:HB3	12:L:54:CYS:O	2.20	0.41
13:M:92:LEU:HD22	13:M:112:ILE:HD11	2.01	0.41
6:T:191:LYS:HB3	6:T:238:TYR:CD1	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:J:1[A]:MET:HE3	10:J:2:GLU:N	2.35	0.41
3:Q:183:THR:HG22	3:Q:186:LEU:HD22	2.02	0.41
9:I:13[A]:MET:HE2	9:I:166:ILE:HB	2.02	0.41
3:Q:239:ASN:HD22	3:Q:239:ASN:N	2.19	0.41
9:W:124:ASP:HB3	15:f:3:ILE:HG23	2.02	0.41
3:C:47:LYS:CB	3:C:48:LYS:CA	2.99	0.41
16:a:302:CL:CL	16:b:302:CL:CL	3.12	0.41
3:C:40:ILE:HD11	3:C:210:VAL:HG13	2.02	0.41
3:C:50:VAL:O	3:C:51:ALA:HB3	2.20	0.41
1:O:10:THR:HG23	20:O:403:HOH:O	2.21	0.41
14:b:48:SER:HB3	14:b:51:ASP:HB2	2.03	0.41
10:J:27[A]:GLN:NE2	20:J:402:HOH:O	2.54	0.41
3:Q:41:VAL:HG11	3:Q:134:VAL:HB	2.02	0.41
4:R:128:ALA:HB1	20:R:424:HOH:O	2.19	0.41
5:S:65[A]:HIS:CE1	20:Z:402:HOH:O	2.74	0.41
10:X:68:LYS:HE3	10:X:69:MET:HE2	2.03	0.41
9:I:124:ASP:OD1	9:I:124:ASP:C	2.64	0.41
7:U:112:ASP:HB3	7:U:152:TYR:CZ	2.56	0.40
9:I:13[A]:MET:HE1	9:I:166:ILE:N	2.35	0.40
11:Y:99:THR:HG22	11:Y:114:VAL:O	2.21	0.40
13:M:86:ARG:NH1	13:M:133:GLU:OE2	2.54	0.40
2:P:151:ASP:HB2	2:P:152:PRO:CD	2.52	0.40
3:C:203:GLY:HA2	3:C:204:LYS:CB	2.51	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	231/234 (99%)	220 (95%)	7 (3%)	4 (2%)	7 9
1	O	228/234 (97%)	217 (95%)	6 (3%)	5 (2%)	5 5

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	248/261 (95%)	236 (95%)	10 (4%)	2 (1%)	16	23
2	P	248/261 (95%)	233 (94%)	13 (5%)	2 (1%)	16	23
3	C	236/248 (95%)	223 (94%)	7 (3%)	6 (2%)	4	4
3	Q	236/248 (95%)	219 (93%)	10 (4%)	7 (3%)	3	3
4	D	232/241 (96%)	223 (96%)	6 (3%)	3 (1%)	9	13
4	R	232/241 (96%)	223 (96%)	6 (3%)	3 (1%)	9	13
5	E	232/263 (88%)	226 (97%)	5 (2%)	1 (0%)	30	42
5	S	238/263 (90%)	231 (97%)	5 (2%)	2 (1%)	16	23
6	F	241/255 (94%)	238 (99%)	3 (1%)	0	100	100
6	T	239/255 (94%)	232 (97%)	5 (2%)	2 (1%)	16	23
7	G	241/246 (98%)	236 (98%)	5 (2%)	0	100	100
7	U	232/246 (94%)	227 (98%)	3 (1%)	2 (1%)	14	20
8	H	220/234 (94%)	216 (98%)	4 (2%)	0	100	100
8	V	220/234 (94%)	217 (99%)	2 (1%)	1 (0%)	24	35
9	I	205/205 (100%)	201 (98%)	3 (2%)	1 (0%)	24	35
9	W	204/205 (100%)	196 (96%)	7 (3%)	1 (0%)	24	35
10	J	195/201 (97%)	191 (98%)	4 (2%)	0	100	100
10	X	195/201 (97%)	193 (99%)	2 (1%)	0	100	100
11	K	198/204 (97%)	195 (98%)	3 (2%)	0	100	100
11	Y	202/204 (99%)	198 (98%)	3 (2%)	1 (0%)	24	35
12	L	213/213 (100%)	211 (99%)	2 (1%)	0	100	100
12	Z	212/213 (100%)	210 (99%)	2 (1%)	0	100	100
13	M	215/219 (98%)	207 (96%)	8 (4%)	0	100	100
13	a	216/219 (99%)	208 (96%)	8 (4%)	0	100	100
14	N	201/205 (98%)	198 (98%)	3 (2%)	0	100	100
14	b	202/205 (98%)	199 (98%)	3 (2%)	0	100	100
15	c	2/5 (40%)	2 (100%)	0	0	100	100
15	d	2/5 (40%)	2 (100%)	0	0	100	100
15	e	2/5 (40%)	2 (100%)	0	0	100	100
15	f	2/5 (40%)	2 (100%)	0	0	100	100
15	g	2/5 (40%)	2 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
15	h	2/5 (40%)	2 (100%)	0	0	100	100
All	All	6224/6488 (96%)	6036 (97%)	145 (2%)	43 (1%)	18	27

All (43) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	52	LYS
1	A	53	SER
3	C	47	LYS
4	D	176	GLY
5	E	59	HIS
1	O	52	LYS
1	O	53	SER
3	Q	47	LYS
3	Q	201	SER
3	Q	206	ILE
4	R	130	PRO
5	S	238	GLU
6	T	208	ALA
11	Y	200	SER
4	D	175[A]	GLU
4	D	175[B]	GLU
1	O	231	ALA
2	P	52	ILE
2	P	54	LYS
3	Q	221	ASN
4	R	128	ALA
7	U	59	LYS
3	C	50	VAL
3	C	200	GLN
3	Q	48	LYS
6	T	6	GLY
8	V	203	ARG
1	A	176	ARG
2	B	51	ASN
2	B	203	VAL
3	C	203	GLY
3	Q	50	VAL
4	R	129	ASP
1	A	201	GLN
3	C	51	ALA
1	O	176	ARG

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Mol	Chain	Res	Type
1	O	201	GLN
5	S	236	LEU
3	C	204	LYS
7	U	58	ASP
9	I	100	GLY
3	Q	203	GLY
9	W	100	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	185/191 (97%)	173 (94%)	12 (6%)	15	25
1	O	176/191 (92%)	167 (95%)	9 (5%)	21	35
2	B	200/221 (90%)	193 (96%)	7 (4%)	32	51
2	P	197/221 (89%)	184 (93%)	13 (7%)	15	25
3	C	179/210 (85%)	171 (96%)	8 (4%)	24	40
3	Q	184/210 (88%)	174 (95%)	10 (5%)	20	33
4	D	189/203 (93%)	183 (97%)	6 (3%)	34	54
4	R	187/203 (92%)	183 (98%)	4 (2%)	47	67
5	E	192/223 (86%)	184 (96%)	8 (4%)	26	43
5	S	197/223 (88%)	191 (97%)	6 (3%)	36	56
6	F	199/212 (94%)	189 (95%)	10 (5%)	22	36
6	T	192/212 (91%)	183 (95%)	9 (5%)	23	39
7	G	202/207 (98%)	196 (97%)	6 (3%)	36	56
7	U	186/207 (90%)	182 (98%)	4 (2%)	45	66
8	H	181/195 (93%)	173 (96%)	8 (4%)	25	41
8	V	172/195 (88%)	160 (93%)	12 (7%)	14	22
9	I	176/174 (101%)	170 (97%)	6 (3%)	32	52
9	W	173/174 (99%)	168 (97%)	5 (3%)	37	57

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
10	J	166/170 (98%)	155 (93%)	11 (7%)	15	25
10	X	163/170 (96%)	155 (95%)	8 (5%)	22	37
11	K	154/159 (97%)	148 (96%)	6 (4%)	28	46
11	Y	159/159 (100%)	156 (98%)	3 (2%)	50	70
12	L	175/178 (98%)	170 (97%)	5 (3%)	37	57
12	Z	175/178 (98%)	171 (98%)	4 (2%)	44	64
13	M	180/181 (99%)	177 (98%)	3 (2%)	53	73
13	a	178/181 (98%)	174 (98%)	4 (2%)	45	66
14	N	158/159 (99%)	155 (98%)	3 (2%)	50	70
14	b	158/159 (99%)	153 (97%)	5 (3%)	34	54
15	c	2/2 (100%)	1 (50%)	1 (50%)	0	0
15	d	2/2 (100%)	1 (50%)	1 (50%)	0	0
15	e	2/2 (100%)	0	2 (100%)	0	0
15	f	2/2 (100%)	1 (50%)	1 (50%)	0	0
15	g	2/2 (100%)	1 (50%)	1 (50%)	0	0
15	h	2/2 (100%)	0	2 (100%)	0	0
All	All	5045/5378 (94%)	4842 (96%)	203 (4%)	28	45

All (203) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	LYS
1	A	61	VAL
1	A	69	LYS
1	A	131	VAL
1	A	135	ILE
1	A	174	GLU
1	A	189	THR
1	A	206	ASN
1	A	221	THR
1	A	223	THR
1	A	226	LYS
1	A	227	ASP
2	B	33	THR
2	B	43	VAL
2	B	59	VAL

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Mol	Chain	Res	Type
2	B	173	SER
2	B	190	LEU
2	B	229	LYS
2	B	249	ARG
3	C	35	VAL
3	C	45	VAL
3	C	148	ASP
3	C	152	THR
3	C	179	GLU
3	C	205	ASN
3	C	208	LEU
3	C	225	ILE
4	D	46	VAL
4	D	126	GLU
4	D	139	VAL
4	D	182	GLN
4	D	199	LEU
4	D	231	LYS
5	E	29	VAL
5	E	61	LYS
5	E	78	THR
5	E	101[A]	ARG
5	E	101[B]	ARG
5	E	181	GLU
5	E	189	LYS
5	E	202	GLU
6	F	17	ASP
6	F	53	VAL
6	F	81	LEU
6	F	86	SER
6	F	87	LEU
6	F	187	ARG
6	F	190	VAL
6	F	240	LYS
6	F	243	LEU
6	F	244	LYS
7	G	42	VAL
7	G	88	ARG
7	G	183	VAL
7	G	192	GLU
7	G	206	LEU
7	G	209	ASP

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Mol	Chain	Res	Type
8	H	6	VAL
8	H	31	CYS
8	H	65	LEU
8	H	68	LEU
8	H	132	LEU
8	H	143	ARG
8	H	215	LYS
8	H	216	ILE
9	I	25[A]	ARG
9	I	25[B]	ARG
9	I	35	THR
9	I	69	ARG
9	I	115	THR
9	I	125	LEU
10	J	1[A]	MET
10	J	1[B]	MET
10	J	27[A]	GLN
10	J	27[B]	GLN
10	J	62	LYS
10	J	88	LEU
10	J	95	ARG
10	J	102	LEU
10	J	155	ARG
10	J	182	ILE
10	J	192	ASP
11	K	12	VAL
11	K	138	VAL
11	K	147	LEU
11	K	158	ARG
11	K	174	VAL
11	K	187	VAL
12	L	3[A]	SER
12	L	3[B]	SER
12	L	73	LYS
12	L	174	LEU
12	L	207	THR
13	M	119	GLU
13	M	154	LEU
13	M	216	SER
14	N	22	THR
14	N	35	THR
14	N	92	GLU

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Mol	Chain	Res	Type
1	O	10	THR
1	O	131	VAL
1	O	135	ILE
1	O	181	LEU
1	O	184	GLU
1	O	189	THR
1	O	206	ASN
1	O	223	THR
1	O	226	LYS
2	P	7[A]	SER
2	P	7[B]	SER
2	P	33	THR
2	P	43	VAL
2	P	53	HIS
2	P	59	VAL
2	P	173	SER
2	P	177	GLN
2	P	190	LEU
2	P	204	SER
2	P	235	GLN
2	P	246	LYS
2	P	249	ARG
3	Q	45	VAL
3	Q	56	GLU
3	Q	98	VAL
3	Q	148	ASP
3	Q	170	GLU
3	Q	206	ILE
3	Q	208	LEU
3	Q	225	ILE
3	Q	227	LYS
3	Q	239	ASN
4	R	46	VAL
4	R	139	VAL
4	R	182	GLN
4	R	209	LYS
5	S	29	VAL
5	S	45	VAL
5	S	54	SER
5	S	101	ARG
5	S	114	SER
5	S	202	GLU

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Mol	Chain	Res	Type
6	T	17	ASP
6	T	53	VAL
6	T	81	LEU
6	T	86	SER
6	T	87	LEU
6	T	190	VAL
6	T	240	LYS
6	T	241	GLU
6	T	243	LEU
7	U	42	VAL
7	U	151	VAL
7	U	199	ILE
7	U	206	LEU
8	V	6	VAL
8	V	31	CYS
8	V	65	LEU
8	V	68	LEU
8	V	104[A]	ASP
8	V	104[B]	ASP
8	V	132	LEU
8	V	143	ARG
8	V	199	LEU
8	V	213	THR
8	V	215	LYS
8	V	216	ILE
9	W	25[A]	ARG
9	W	25[B]	ARG
9	W	35	THR
9	W	69	ARG
9	W	125	LEU
10	X	1	MET
10	X	62	LYS
10	X	88	LEU
10	X	95	ARG
10	X	102	LEU
10	X	132	HIS
10	X	158	GLU
10	X	182	ILE
11	Y	12	VAL
11	Y	138	VAL
11	Y	147	LEU
12	Z	142	SER

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Mol	Chain	Res	Type
12	Z	174	LEU
12	Z	207	THR
12	Z	208	VAL
13	a	92	LEU
13	a	119	GLU
13	a	154	LEU
13	a	216	SER
14	b	22	THR
14	b	35	THR
14	b	92	GLU
14	b	145	GLU
14	b	202	LEU
15	c	3	ILE
15	d	3	ILE
15	e	3	ILE
15	e	4	THR
15	f	3	ILE
15	g	3	ILE
15	h	3	ILE
15	h	4	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (67) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	51	GLN
1	A	101	GLN
1	A	108	GLN
2	B	40	ASN
2	B	51	ASN
3	C	54	GLN
3	C	94	HIS
4	D	182	GLN
5	E	8	ASN
5	E	60	GLN
5	E	65	HIS
6	F	143	ASN
7	G	12	HIS
7	G	68	HIS
7	G	75	ASN
7	G	128	ASN
8	H	80	ASN
8	H	172	ASN

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Mol	Chain	Res	Type
8	H	193	ASN
10	J	61	GLN
10	J	63	ASN
10	J	101	ASN
10	J	132	HIS
11	K	62	GLN
12	L	163	HIS
13	M	65	GLN
13	M	162	GLN
13	M	188	GLN
14	N	106	GLN
1	O	101	GLN
1	O	118	GLN
2	P	40	ASN
2	P	109	GLN
2	P	142	HIS
2	P	146	GLN
2	P	166	ASN
3	Q	18	GLN
3	Q	94	HIS
3	Q	239	ASN
4	R	152	GLN
4	R	182	GLN
4	R	186	HIS
5	S	8	ASN
5	S	60	GLN
5	S	86	ASN
6	T	63	ASN
6	T	68	ASN
6	T	143	ASN
7	U	68	HIS
7	U	75	ASN
7	U	90	GLN
7	U	128	ASN
8	V	80	ASN
8	V	172	ASN
8	V	193	ASN
10	X	24	ASN
10	X	61	GLN
10	X	63	ASN
10	X	174	ASN
11	Y	62	GLN

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Mol	Chain	Res	Type
12	Z	58	HIS
12	Z	79	ASN
12	Z	157	ASN
13	a	65	GLN
13	a	162	GLN
13	a	188	GLN
14	b	193	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
7	YCM	U	137	7	7,9,10	0.87	0	5,10,12	0.62	0
10	6V1	X	91	10	13,15,16	1.87	3 (23%)	10,20,22	3.87	6 (60%)
3	YCM	Q	63	3	7,9,10	1.11	1 (14%)	5,10,12	2.36	2 (40%)
15	IML	h	2	15	7,8,9	0.45	0	6,9,11	1.54	1 (16%)
7	6V1	U	47	7	13,15,16	1.87	4 (30%)	10,20,22	2.41	3 (30%)
7	6V1	G	47	7	13,15,16	2.23	4 (30%)	10,20,22	2.15	3 (30%)
15	IML	d	2	15	7,8,9	0.98	0	6,9,11	2.52	3 (50%)
7	6V1	U	161	7	13,15,16	1.94	4 (30%)	10,20,22	2.29	4 (40%)
15	IML	c	2	15	7,8,9	0.97	0	6,9,11	1.69	3 (50%)
3	YCM	C	63	3	7,9,10	0.97	1 (14%)	5,10,12	0.70	0
15	IML	f	2	15	7,8,9	0.73	0	6,9,11	1.55	1 (16%)
5	6V1	S	148	5	13,15,16	2.14	4 (30%)	10,20,22	2.76	5 (50%)
7	6V1	G	161	7	13,15,16	2.11	4 (30%)	10,20,22	2.11	4 (40%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
10	6V1	J	91	10	13,15,16	1.81	3 (23%)	10,20,22	4.02	6 (60%)
15	IML	e	2	15	7,8,9	0.76	0	6,9,11	2.59	1 (16%)
5	6V1	E	148	5	13,15,16	2.27	4 (30%)	10,20,22	3.16	5 (50%)
15	IML	g	2	15	7,8,9	1.00	1 (14%)	6,9,11	2.37	2 (33%)
7	YCM	G	137	7	7,9,10	0.99	0	5,10,12	1.34	1 (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
7	YCM	U	137	7	-	2/6/8/10	-
10	6V1	X	91	10	-	4/6/25/27	0/1/1/1
3	YCM	Q	63	3	-	5/6/8/10	-
15	IML	h	2	15	2/2/2/4	1/8/10/12	-
7	6V1	U	47	7	1/1/5/6	1/6/25/27	0/1/1/1
15	IML	d	2	15	2/2/2/4	2/8/10/12	-
15	IML	f	2	15	1/1/2/4	5/8/10/12	-
15	IML	c	2	15	1/1/2/4	3/8/10/12	-
15	IML	g	2	15	2/2/2/4	1/8/10/12	-
3	YCM	C	63	3	-	1/6/8/10	-
7	6V1	G	47	7	-	2/6/25/27	0/1/1/1
5	6V1	S	148	5	-	2/6/25/27	0/1/1/1
7	6V1	G	161	7	-	1/6/25/27	0/1/1/1
15	IML	e	2	15	1/1/2/4	6/8/10/12	-
7	6V1	U	161	7	-	1/6/25/27	0/1/1/1
5	6V1	E	148	5	-	2/6/25/27	0/1/1/1
10	6V1	J	91	10	-	4/6/25/27	0/1/1/1
7	YCM	G	137	7	-	1/6/8/10	-

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	S	148	6V1	CB-SG	-5.96	1.76	1.82
5	E	148	6V1	CB-SG	-5.68	1.76	1.82
7	G	161	6V1	CB-SG	-5.60	1.76	1.82
7	G	47	6V1	CB-SG	-5.35	1.76	1.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	U	47	6V1	CB-SG	-4.79	1.77	1.82
7	U	161	6V1	CB-SG	-4.59	1.77	1.82
10	X	91	6V1	CB-SG	-4.43	1.77	1.82
5	E	148	6V1	C1-SG	-4.19	1.78	1.83
10	J	91	6V1	C1-SG	-4.12	1.78	1.83
7	G	47	6V1	C4-N3	-3.96	1.32	1.38
7	U	161	6V1	C1-SG	-3.42	1.79	1.83
7	G	47	6V1	C2-N3	-3.36	1.34	1.38
10	X	91	6V1	C1-SG	-3.36	1.79	1.83
10	J	91	6V1	C4-N3	-3.19	1.33	1.38
10	X	91	6V1	C4-N3	-3.10	1.34	1.38
10	J	91	6V1	CB-SG	-3.10	1.79	1.82
7	G	161	6V1	C2-N3	-3.09	1.34	1.38
5	S	148	6V1	C1-SG	-3.00	1.80	1.83
5	E	148	6V1	C2-N3	-2.98	1.34	1.38
7	G	161	6V1	C1-SG	-2.75	1.80	1.83
5	S	148	6V1	C2-N3	-2.71	1.35	1.38
7	U	161	6V1	C4-N3	-2.59	1.34	1.38
7	G	161	6V1	C4-N3	-2.53	1.34	1.38
3	Q	63	YCM	CD-SG	-2.53	1.75	1.81
7	U	47	6V1	C2-N3	-2.52	1.35	1.38
7	U	161	6V1	C2-N3	-2.49	1.35	1.38
7	U	47	6V1	C1-SG	-2.46	1.80	1.83
5	E	148	6V1	C4-N3	-2.38	1.35	1.38
7	G	47	6V1	C1-SG	-2.33	1.80	1.83
7	U	47	6V1	C4-N3	-2.33	1.35	1.38
15	g	2	IML	CA-N	-2.25	1.43	1.47
3	C	63	YCM	CD-SG	-2.09	1.76	1.81
5	S	148	6V1	C4-N3	-2.02	1.35	1.38

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
10	J	91	6V1	C5-C4-N3	7.29	112.66	108.07
10	X	91	6V1	C5-C4-N3	6.74	112.31	108.07
10	J	91	6V1	C2-N3-C4	-6.34	109.36	113.07
10	X	91	6V1	C2-N3-C4	-6.26	109.41	113.07
5	E	148	6V1	C5-C4-N3	6.19	111.97	108.07
15	e	2	IML	CB-CA-C	-6.06	104.47	112.77
5	E	148	6V1	C2-N3-C4	-5.89	109.62	113.07
7	U	47	6V1	C2-N3-C4	-5.49	109.85	113.07
10	J	91	6V1	C6-N3-C2	5.14	129.40	123.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	S	148	6V1	C5-C4-N3	4.97	111.20	108.07
5	S	148	6V1	C2-N3-C4	-4.96	110.16	113.07
10	X	91	6V1	C6-N3-C2	4.80	129.00	123.37
10	X	91	6V1	O7-C2-N3	4.56	129.67	124.14
10	J	91	6V1	O7-C2-N3	4.54	129.65	124.14
7	G	47	6V1	C2-N3-C4	-4.38	110.50	113.07
15	g	2	IML	CG2-CB-CA	4.33	121.27	110.99
3	Q	63	YCM	CE-CD-SG	-4.19	100.59	113.81
7	U	47	6V1	C5-C4-N3	3.86	110.50	108.07
7	U	161	6V1	C5-C4-N3	3.77	110.44	108.07
15	d	2	IML	CG2-CB-CA	3.70	119.79	110.99
7	U	161	6V1	C2-N3-C4	-3.53	111.00	113.07
15	d	2	IML	CB-CA-C	3.43	117.47	112.77
7	U	161	6V1	O8-C4-C5	-3.41	122.29	127.28
7	G	161	6V1	O8-C4-N3	3.39	127.57	123.94
7	G	161	6V1	O8-C4-C5	-3.35	122.37	127.28
10	J	91	6V1	O8-C4-C5	-3.28	122.48	127.28
10	X	91	6V1	O8-C4-C5	-3.25	122.53	127.28
5	E	148	6V1	O8-C4-C5	-3.20	122.59	127.28
7	G	47	6V1	C5-C4-N3	3.19	110.08	108.07
7	G	161	6V1	C5-C4-N3	3.15	110.05	108.07
7	U	161	6V1	O8-C4-N3	3.10	127.26	123.94
5	S	148	6V1	O8-C4-C5	-3.01	122.88	127.28
15	d	2	IML	CD1-CG1-CB	2.91	125.90	113.81
7	G	161	6V1	C2-N3-C4	-2.80	111.43	113.07
15	h	2	IML	O-C-CA	-2.78	117.51	124.86
10	X	91	6V1	O7-C2-C1	-2.76	118.50	124.52
7	G	137	YCM	CE-CD-SG	2.74	122.46	113.81
5	E	148	6V1	C6-N3-C4	2.62	125.62	122.52
5	S	148	6V1	CB-CA-C	-2.59	103.75	110.80
7	U	47	6V1	C6-N3-C2	2.54	126.35	123.37
10	J	91	6V1	O7-C2-C1	-2.54	118.99	124.52
15	g	2	IML	CG1-CB-CA	-2.41	105.12	111.17
5	S	148	6V1	C6-N3-C4	2.39	125.35	122.52
15	c	2	IML	CG2-CB-CA	2.39	116.67	110.99
5	E	148	6V1	CB-CA-C	-2.39	104.32	110.80
3	Q	63	YCM	CA-CB-SG	-2.34	104.88	113.51
7	G	47	6V1	O8-C4-C5	-2.34	123.86	127.28
15	f	2	IML	O-C-CA	-2.20	119.04	124.86
15	c	2	IML	CG1-CB-CA	-2.12	105.86	111.17
15	c	2	IML	O-C-CA	-2.08	119.37	124.86

All (10) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
7	U	47	6V1	C1
15	c	2	IML	CB
15	d	2	IML	CB
15	d	2	IML	CA
15	e	2	IML	CA
15	f	2	IML	CB
15	g	2	IML	CB
15	g	2	IML	CA
15	h	2	IML	CB
15	h	2	IML	CA

All (44) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	Q	63	YCM	SG-CD-CE-OZ1
3	Q	63	YCM	SG-CD-CE-NZ2
7	U	137	YCM	SG-CD-CE-NZ2
5	E	148	6V1	C3-C6-N3-C2
5	S	148	6V1	C3-C6-N3-C2
5	S	148	6V1	C3-C6-N3-C4
7	G	47	6V1	C-CA-CB-SG
7	U	47	6V1	C-CA-CB-SG
10	J	91	6V1	C-CA-CB-SG
10	J	91	6V1	C3-C6-N3-C2
10	J	91	6V1	C3-C6-N3-C4
10	X	91	6V1	C-CA-CB-SG
10	X	91	6V1	C3-C6-N3-C2
10	X	91	6V1	C3-C6-N3-C4
15	d	2	IML	C-CA-CB-CG2
15	e	2	IML	C-CA-CB-CG2
15	e	2	IML	C-CA-CB-CG1
15	f	2	IML	O-C-CA-CB
15	f	2	IML	N-CA-CB-CG2
15	f	2	IML	C-CA-CB-CG2
15	g	2	IML	CB-CA-N-CN
15	h	2	IML	CB-CA-N-CN
15	e	2	IML	CG2-CB-CG1-CD1
15	f	2	IML	N-CA-CB-CG1
5	E	148	6V1	C3-C6-N3-C4
15	e	2	IML	CA-CB-CG1-CD1
15	c	2	IML	CG2-CB-CG1-CD1
15	c	2	IML	CA-CB-CG1-CD1
3	Q	63	YCM	N-CA-CB-SG

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Mol	Chain	Res	Type	Atoms
3	C	63	YCM	CE-CD-SG-CB
7	G	137	YCM	CE-CD-SG-CB
3	Q	63	YCM	CE-CD-SG-CB
7	U	137	YCM	CE-CD-SG-CB
15	f	2	IML	CG2-CB-CG1-CD1
10	J	91	6V1	CA-CB-SG-C1
10	X	91	6V1	CA-CB-SG-C1
7	G	47	6V1	C5-C1-SG-CB
15	c	2	IML	CB-CA-N-CN
15	e	2	IML	CB-CA-N-CN
15	d	2	IML	N-CA-CB-CG2
15	e	2	IML	N-CA-CB-CG1
7	G	161	6V1	N-CA-CB-SG
7	U	161	6V1	N-CA-CB-SG
3	Q	63	YCM	CA-CB-SG-CD

There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
15	d	2	IML	3	0
15	c	2	IML	1	0
15	f	2	IML	1	0
15	g	2	IML	3	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 83 ligands modelled in this entry, 74 are monoatomic - leaving 9 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
19	1PE	I	303	-	15,15,15	0.58	0	14,14,14	0.87	1 (7%)
19	1PE	N	305	-	15,15,15	0.53	0	14,14,14	0.32	0
19	1PE	H	304	-	15,15,15	0.57	0	14,14,14	0.51	0
19	1PE	Z	301	-	15,15,15	0.57	0	14,14,14	0.42	0
19	1PE	L	301	-	15,15,15	0.55	0	14,14,14	0.33	0
19	1PE	a	304	-	15,15,15	0.56	0	14,14,14	0.25	0
19	1PE	W	303	-	15,15,15	0.63	0	14,14,14	0.48	0
19	1PE	K	306	-	15,15,15	0.57	0	14,14,14	0.61	0
19	1PE	b	305	-	15,15,15	0.61	0	14,14,14	0.47	0

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
19	1PE	I	303	-	-	7/13/13/13	-
19	1PE	N	305	-	-	5/13/13/13	-
19	1PE	H	304	-	-	5/13/13/13	-
19	1PE	Z	301	-	-	7/13/13/13	-
19	1PE	L	301	-	-	3/13/13/13	-
19	1PE	a	304	-	-	7/13/13/13	-
19	1PE	W	303	-	-	10/13/13/13	-
19	1PE	K	306	-	-	9/13/13/13	-
19	1PE	b	305	-	-	6/13/13/13	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
19	I	303	1PE	C25-OH5-C14	2.16	122.72	113.26

There are no chirality outliers.

All (59) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
19	I	303	1PE	C15-C25-OH5-C14
19	H	304	1PE	C24-C14-OH5-C25
19	K	306	1PE	C16-C26-OH6-C15

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Mol	Chain	Res	Type	Atoms
19	b	305	1PE	OH4-C13-C23-OH3
19	W	303	1PE	OH6-C15-C25-OH5
19	Z	301	1PE	OH6-C15-C25-OH5
19	H	304	1PE	OH4-C13-C23-OH3
19	K	306	1PE	OH5-C14-C24-OH4
19	a	304	1PE	OH5-C14-C24-OH4
19	b	305	1PE	OH6-C15-C25-OH5
19	a	304	1PE	OH4-C13-C23-OH3
19	W	303	1PE	OH7-C16-C26-OH6
19	N	305	1PE	OH2-C12-C22-OH3
19	L	301	1PE	OH6-C15-C25-OH5
19	N	305	1PE	OH7-C16-C26-OH6
19	a	304	1PE	OH2-C12-C22-OH3
19	Z	301	1PE	OH4-C13-C23-OH3
19	N	305	1PE	OH4-C13-C23-OH3
19	L	301	1PE	OH5-C14-C24-OH4
19	K	306	1PE	OH6-C15-C25-OH5
19	I	303	1PE	OH6-C15-C25-OH5
19	a	304	1PE	OH7-C16-C26-OH6
19	b	305	1PE	OH7-C16-C26-OH6
19	Z	301	1PE	C16-C26-OH6-C15
19	I	303	1PE	OH2-C12-C22-OH3
19	H	304	1PE	OH5-C14-C24-OH4
19	Z	301	1PE	C15-C25-OH5-C14
19	H	304	1PE	C16-C26-OH6-C15
19	W	303	1PE	C14-C24-OH4-C13
19	W	303	1PE	C24-C14-OH5-C25
19	I	303	1PE	C14-C24-OH4-C13
19	Z	301	1PE	C12-C22-OH3-C23
19	W	303	1PE	C25-C15-OH6-C26
19	I	303	1PE	C24-C14-OH5-C25
19	W	303	1PE	C12-C22-OH3-C23
19	K	306	1PE	C13-C23-OH3-C22
19	Z	301	1PE	OH7-C16-C26-OH6
19	a	304	1PE	C23-C13-OH4-C24
19	Z	301	1PE	C25-C15-OH6-C26
19	a	304	1PE	OH6-C15-C25-OH5
19	K	306	1PE	OH4-C13-C23-OH3
19	W	303	1PE	C13-C23-OH3-C22
19	K	306	1PE	C25-C15-OH6-C26
19	N	305	1PE	C12-C22-OH3-C23
19	H	304	1PE	OH6-C15-C25-OH5

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Mol	Chain	Res	Type	Atoms
19	L	301	1PE	OH4-C13-C23-OH3
19	I	303	1PE	C13-C23-OH3-C22
19	a	304	1PE	C25-C15-OH6-C26
19	b	305	1PE	C24-C14-OH5-C25
19	W	303	1PE	C23-C13-OH4-C24
19	K	306	1PE	OH2-C12-C22-OH3
19	W	303	1PE	OH5-C14-C24-OH4
19	b	305	1PE	C14-C24-OH4-C13
19	b	305	1PE	OH5-C14-C24-OH4
19	K	306	1PE	C14-C24-OH4-C13
19	N	305	1PE	C13-C23-OH3-C22
19	W	303	1PE	C16-C26-OH6-C15
19	I	303	1PE	OH5-C14-C24-OH4
19	K	306	1PE	C24-C14-OH5-C25

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	230/234 (98%)	-0.36	1 (0%) 88 87	31, 60, 99, 113	3 (1%)
1	O	230/234 (98%)	-0.07	0 100 100	56, 83, 125, 146	0
2	B	248/261 (95%)	-0.32	0 100 100	30, 67, 117, 159	2 (0%)
2	P	248/261 (95%)	-0.20	2 (0%) 82 80	42, 75, 126, 169	2 (0%)
3	C	236/248 (95%)	0.02	0 100 100	32, 79, 125, 152	2 (0%)
3	Q	238/248 (95%)	0.10	6 (2%) 58 55	49, 80, 144, 184	0
4	D	233/241 (96%)	-0.12	2 (0%) 81 79	37, 72, 105, 135	1 (0%)
4	R	233/241 (96%)	-0.27	3 (1%) 75 72	29, 59, 94, 124	1 (0%)
5	E	233/263 (88%)	-0.39	1 (0%) 88 87	31, 56, 104, 129	1 (0%)
5	S	237/263 (90%)	-0.26	3 (1%) 75 72	31, 65, 109, 124	3 (1%)
6	F	239/255 (93%)	-0.53	1 (0%) 88 87	28, 48, 72, 93	4 (1%)
6	T	240/255 (94%)	0.04	4 (1%) 69 66	43, 76, 112, 144	1 (0%)
7	G	241/246 (97%)	-0.49	0 100 100	30, 54, 93, 131	2 (0%)
7	U	235/246 (95%)	-0.03	4 (1%) 69 66	44, 85, 119, 162	1 (0%)
8	H	220/234 (94%)	-0.59	0 100 100	28, 51, 85, 126	2 (0%)
8	V	220/234 (94%)	-0.26	2 (0%) 81 79	31, 64, 100, 125	2 (0%)
9	I	204/205 (99%)	-0.58	0 100 100	29, 51, 75, 89	3 (1%)
9	W	204/205 (99%)	-0.37	1 (0%) 87 86	32, 63, 93, 102	2 (0%)
10	J	195/201 (97%)	-0.49	1 (0%) 87 86	22, 56, 78, 94	3 (1%)
10	X	195/201 (97%)	-0.44	0 100 100	25, 57, 74, 92	2 (1%)
11	K	200/204 (98%)	-0.47	1 (0%) 87 86	44, 60, 87, 100	0
11	Y	201/204 (98%)	-0.57	2 (0%) 79 77	29, 51, 77, 95	3 (1%)
12	L	213/213 (100%)	-0.46	0 100 100	32, 60, 86, 108	2 (0%)
12	Z	213/213 (100%)	-0.59	0 100 100	36, 52, 79, 96	1 (0%)

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
13	M	216/219 (98%)	-0.49	1 (0%) 87 86	38, 53, 81, 112	1 (0%)
13	a	216/219 (98%)	-0.48	0 100 100	35, 55, 78, 101	2 (0%)
14	N	202/205 (98%)	-0.60	1 (0%) 87 86	28, 49, 77, 109	1 (0%)
14	b	203/205 (99%)	-0.44	0 100 100	43, 59, 89, 127	1 (0%)
15	c	2/5 (40%)	0.28	0 100 100	49, 49, 49, 55	0
15	d	2/5 (40%)	-0.02	0 100 100	51, 51, 51, 56	0
15	e	2/5 (40%)	-0.21	0 100 100	51, 51, 51, 51	0
15	f	2/5 (40%)	0.20	0 100 100	62, 62, 62, 63	0
15	g	2/5 (40%)	-0.21	0 100 100	42, 42, 42, 49	0
15	h	2/5 (40%)	0.11	0 100 100	59, 59, 59, 64	0
All	All	6235/6488 (96%)	-0.34	36 (0%) 85 84	22, 61, 108, 184	48 (0%)

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
6	T	5	THR	4.1
11	Y	201	GLY	3.7
4	R	241	ILE	3.7
5	S	18[A]	ARG	3.5
8	V	204	CYS	3.5
6	T	208	ALA	3.1
7	U	193	GLN	3.0
11	K	40	TYR	2.9
7	U	186	LYS	2.8
7	U	185	LYS	2.8
5	S	2	PHE	2.8
3	Q	50	VAL	2.7
11	Y	40	TYR	2.6
3	Q	48	LYS	2.6
14	N	202	LEU	2.6
9	W	16[A]	LYS	2.6
7	U	245	ARG	2.5
8	V	220	GLU	2.5
2	P	61	PHE	2.4
5	S	239	ARG	2.4
4	R	131	GLY	2.3
4	D	241	ILE	2.3
5	E	56	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	3	ARG	2.3
6	T	6	GLY	2.3
3	Q	202	GLY	2.3
3	Q	51	ALA	2.2
4	R	120[A]	ALA	2.2
6	F	244	LYS	2.2
3	Q	85	ASN	2.2
13	M	215	ILE	2.2
2	P	203	VAL	2.1
6	T	209	PHE	2.1
4	D	131	GLY	2.1
3	Q	232	ILE	2.1
10	J	1[A]	MET	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
7	6V1	U	47	15/16	0.87	0.15	103,127,131,132	0
7	YCM	G	137	10/11	0.89	0.11	48,56,70,71	0
15	IML	d	2	9/10	0.89	0.13	54,56,59,60	0
7	6V1	U	161	15/16	0.90	0.11	78,92,98,101	0
7	YCM	U	137	10/11	0.91	0.11	75,84,96,96	0
3	YCM	Q	63	10/11	0.91	0.09	67,70,76,77	0
15	IML	g	2	9/10	0.91	0.13	55,57,60,64	0
3	YCM	C	63	10/11	0.92	0.09	69,74,80,82	0
15	IML	f	2	9/10	0.92	0.11	64,71,73,74	0
5	6V1	S	148	15/16	0.92	0.11	56,86,92,92	0
15	IML	e	2	9/10	0.93	0.12	58,62,68,69	0
15	IML	h	2	9/10	0.93	0.12	67,68,75,76	0
15	IML	c	2	9/10	0.94	0.10	52,59,59,62	0
5	6V1	E	148	15/16	0.94	0.10	44,70,77,81	0
7	6V1	G	47	15/16	0.94	0.10	54,68,71,72	0
7	6V1	G	161	15/16	0.95	0.09	47,71,80,84	0
10	6V1	X	91	15/16	0.95	0.10	55,73,77,77	0
10	6V1	J	91	15/16	0.96	0.10	50,71,74,78	0

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
16	CL	O	304	1/1	0.85	0.14	91,91,91,91	0
19	1PE	H	304	16/16	0.87	0.16	85,100,110,110	0
16	CL	Y	305	1/1	0.88	0.12	84,84,84,84	0
16	CL	D	301	1/1	0.88	0.16	90,90,90,90	0
19	1PE	W	303	16/16	0.88	0.12	78,85,94,95	0
16	CL	S	302	1/1	0.89	0.12	95,95,95,95	0
16	CL	E	303	1/1	0.89	0.14	91,91,91,91	0
16	CL	a	303	1/1	0.89	0.11	84,84,84,84	0
16	CL	Q	301	1/1	0.89	0.08	97,97,97,97	0
19	1PE	L	301	16/16	0.89	0.17	95,101,126,127	0
16	CL	S	301	1/1	0.89	0.22	95,95,95,95	0
19	1PE	b	305	16/16	0.89	0.16	67,72,100,102	0
19	1PE	I	303	16/16	0.90	0.12	68,81,97,102	0
19	1PE	Z	301	16/16	0.90	0.13	74,89,98,101	0
19	1PE	a	304	16/16	0.90	0.15	82,88,121,123	0
16	CL	Q	302	1/1	0.90	0.16	89,89,89,89	0
16	CL	Y	303	1/1	0.91	0.12	89,89,89,89	0
16	CL	S	303	1/1	0.91	0.17	82,82,82,82	0
19	1PE	K	306	16/16	0.91	0.12	75,87,95,97	0
16	CL	V	303	1/1	0.91	0.16	75,75,75,75	0
16	CL	O	303	1/1	0.92	0.08	101,101,101,101	0
16	CL	C	302	1/1	0.92	0.17	86,86,86,86	0
16	CL	K	305	1/1	0.92	0.15	80,80,80,80	0
16	CL	N	303	1/1	0.92	0.13	87,87,87,87	0
16	CL	R	301	1/1	0.92	0.10	84,84,84,84	0
16	CL	Y	306	1/1	0.92	0.20	90,90,90,90	0
16	CL	O	302	1/1	0.92	0.11	79,79,79,79	0
17	K	b	306	1/1	0.92	0.08	73,73,73,73	0
17	K	L	302	1/1	0.93	0.08	78,78,78,78	0
16	CL	K	304	1/1	0.93	0.10	89,89,89,89	0
16	CL	C	301	1/1	0.93	0.11	82,82,82,82	0
16	CL	A	304	1/1	0.93	0.12	77,77,77,77	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
16	CL	R	302	1/1	0.93	0.32	81,81,81,81	0
16	CL	N	304	1/1	0.94	0.14	72,72,72,72	0
16	CL	b	302	1/1	0.94	0.12	91,91,91,91	0
16	CL	b	303	1/1	0.94	0.09	82,82,82,82	0
19	1PE	N	305	16/16	0.94	0.10	55,61,82,83	0
16	CL	O	301	1/1	0.94	0.10	80,80,80,80	0
17	K	N	306	1/1	0.94	0.08	65,65,65,65	0
16	CL	E	304	1/1	0.94	0.10	81,81,81,81	0
16	CL	A	303	1/1	0.94	0.09	72,72,72,72	0
16	CL	U	301	1/1	0.95	0.14	79,79,79,79	0
17	K	Z	302	1/1	0.95	0.06	63,63,63,63	0
16	CL	B	302	1/1	0.95	0.15	80,80,80,80	0
16	CL	W	302	1/1	0.95	0.07	70,70,70,70	0
16	CL	E	302	1/1	0.95	0.16	79,79,79,79	0
16	CL	I	302	1/1	0.95	0.08	64,64,64,64	0
16	CL	G	302	1/1	0.96	0.05	79,79,79,79	0
16	CL	A	301	1/1	0.96	0.07	64,64,64,64	0
16	CL	Y	304	1/1	0.96	0.07	77,77,77,77	0
16	CL	K	303	1/1	0.96	0.07	84,84,84,84	0
16	CL	A	302	1/1	0.96	0.06	80,80,80,80	0
16	CL	a	301	1/1	0.96	0.09	88,88,88,88	0
16	CL	E	301	1/1	0.96	0.12	85,85,85,85	0
16	CL	M	301	1/1	0.96	0.16	86,86,86,86	0
16	CL	P	301	1/1	0.96	0.10	67,67,67,67	0
16	CL	c	101	1/1	0.96	0.14	63,63,63,63	0
17	K	G	303	1/1	0.96	0.07	56,56,56,56	0
16	CL	M	303	1/1	0.96	0.06	72,72,72,72	0
18	MG	I	301	1/1	0.97	0.06	60,60,60,60	0
18	MG	W	301	1/1	0.97	0.08	56,56,56,56	0
16	CL	M	302	1/1	0.97	0.10	59,59,59,59	0
16	CL	f	101	1/1	0.97	0.07	64,64,64,64	0
16	CL	a	302	1/1	0.97	0.10	64,64,64,64	0
16	CL	H	303	1/1	0.97	0.12	70,70,70,70	0
16	CL	b	301	1/1	0.97	0.08	62,62,62,62	0
17	K	U	302	1/1	0.97	0.06	72,72,72,72	0
16	CL	N	302	1/1	0.97	0.07	53,53,53,53	0
16	CL	F	301	1/1	0.97	0.09	79,79,79,79	0
18	MG	H	302	1/1	0.97	0.13	63,63,63,63	0
16	CL	G	301	1/1	0.98	0.25	65,65,65,65	0
18	MG	H	301	1/1	0.98	0.03	64,64,64,64	0
16	CL	B	301	1/1	0.98	0.09	53,53,53,53	0
16	CL	b	304	1/1	0.98	0.06	55,55,55,55	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
18	MG	K	301	1/1	0.98	0.06	52,52,52,52	0
16	CL	K	302	1/1	0.98	0.10	54,54,54,54	0
18	MG	Y	301	1/1	0.98	0.07	47,47,47,47	0
16	CL	Y	302	1/1	0.98	0.06	49,49,49,49	0
18	MG	V	302	1/1	0.99	0.03	72,72,72,72	0
18	MG	J	301	1/1	0.99	0.03	66,66,66,66	0
16	CL	N	301	1/1	0.99	0.04	47,47,47,47	0
18	MG	V	301	1/1	0.99	0.08	55,55,55,55	0
18	MG	X	301	1/1	1.00	0.04	54,54,54,54	0

6.5 Other polymers [i](#)

There are no such residues in this entry.