



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 08:17 PM UTC

PDB ID : 4Y8S / pdb_00004y8s
Title : Yeast 20S proteasome beta2-H116D mutant in complex with Ac-LAE-ep
Authors : Huber, E.M.; Groll, M.
Deposited on : 2015-02-16
Resolution : 2.70 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

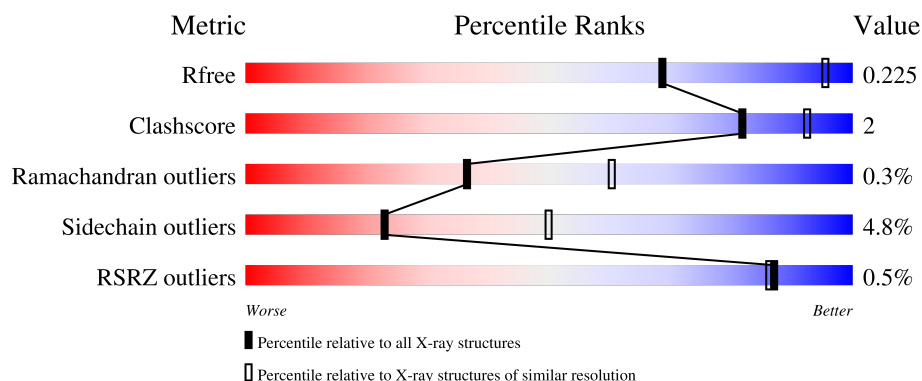
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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

Mol	Chain	Length	Quality of chain
1	A	250	96%
1	O	250	96%
2	B	258	82% 11% 5%
2	P	258	83% 10% 5%
3	C	254	84% 8% 6%

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Mol	Chain	Length	Quality of chain
3	Q	254	 2% 84% 8% 6%
4	D	260	 80% 10% 10%
4	R	260	 78% 11% 10%
5	E	234	 90% 9%
5	S	234	 89% 9%
6	F	288	 80% 16%
6	T	288	 2% 80% 16%
7	G	252	 89% 5%
7	U	252	 88% 6%
8	H	232	 88% 7%
8	V	232	 90% 5%
9	I	205	 90% 9%
9	W	205	 90% 9%
10	J	198	 2% 87% 10%
10	X	198	 2% 87% 10%
11	K	212	 92% 8%
11	Y	212	 91% 8%
12	L	222	 91% 8%
12	Z	222	 89% 10%
13	M	246	 84% 8% 5%
13	a	246	 84% 9% 5%
14	N	196	 92% 7%
14	b	196	 92% 7%
15	c	5	 60% 40%
15	d	5	 80% 20%

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Mol	Chain	Length	Quality of chain
15	e	5	20%60%20%
15	f	5	80%20%
15	g	5	60%20%20%
15	h	5	60%40%

2 Entry composition

There are 18 unique types of molecules in this entry. The entry contains 49825 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
			1682	1059	291	325	7			
8	V	222	Total	C	N	O	S	0	1	0
			1689	1064	293	325	7			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
H	116	ASP	HIS	engineered mutation	UNP P25043
V	116	ASP	HIS	engineered mutation	UNP P25043

- 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			
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	1	0
			1764	1120	305	335	4			

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

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

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

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

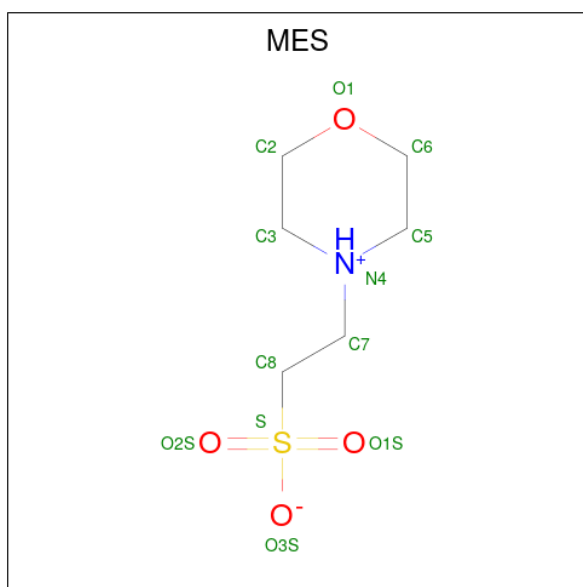
- Molecule 15 is a protein called Ac-LAE-ep.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
15	c	5	Total	C	N	O	0	0	0
			29	19	3	7			
15	d	5	Total	C	N	O	0	0	0
			29	19	3	7			
15	e	5	Total	C	N	O	0	0	0
			29	19	3	7			
15	f	5	Total	C	N	O	0	0	0
			29	19	3	7			
15	g	5	Total	C	N	O	0	0	0
			29	19	3	7			
15	h	5	Total	C	N	O	0	0	0
			29	19	3	7			

- Molecule 16 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 17 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (CCD ID: MES) (formula: C₆H₁₃NO₄S).



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

- Molecule 18 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
18	A	6	Total	O	0	0
			6	6		
18	B	13	Total	O	0	0
			13	13		
18	C	10	Total	O	0	0
			10	10		
18	D	2	Total	O	0	0
			2	2		
18	E	5	Total	O	0	0
			5	5		
18	F	13	Total	O	0	0
			13	13		
18	G	15	Total	O	0	0
			15	15		
18	H	23	Total	O	0	0
			23	23		
18	I	10	Total	O	0	0
			10	10		
18	J	11	Total	O	0	0
			11	11		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
18	K	14	Total O 14 14	0	0
18	L	14	Total O 14 14	0	0
18	M	13	Total O 13 13	0	0
18	N	6	Total O 6 6	0	0
18	O	3	Total O 3 3	0	0
18	P	7	Total O 7 7	0	0
18	Q	8	Total O 8 8	0	0
18	R	3	Total O 3 3	0	0
18	S	4	Total O 4 4	0	0
18	T	9	Total O 9 9	0	0
18	U	21	Total O 21 21	0	0
18	V	13	Total O 13 13	0	0
18	W	8	Total O 8 8	0	0
18	X	14	Total O 14 14	0	0
18	Y	16	Total O 16 16	0	0
18	Z	20	Total O 20 20	0	0
18	a	17	Total O 17 17	0	0
18	b	11	Total O 11 11	0	0
18	c	1	Total O 1 1	0	0
18	d	1	Total O 1 1	0	0
18	e	1	Total O 1 1	0	0

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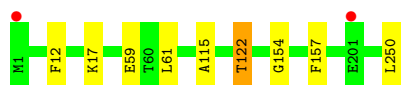
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
18	f	1	Total	O	0	0
			1	1		
18	g	1	Total	O	0	0
			1	1		

3 Residue-property plots [i](#)

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

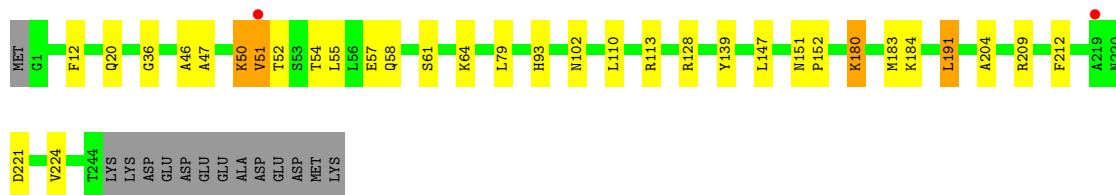
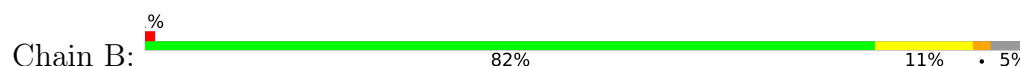
- Molecule 1: Proteasome subunit alpha type-2



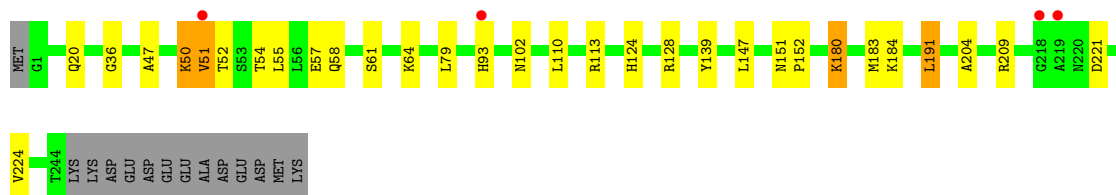
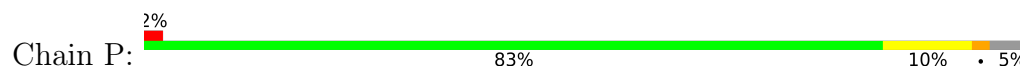
- Molecule 1: Proteasome subunit alpha type-2



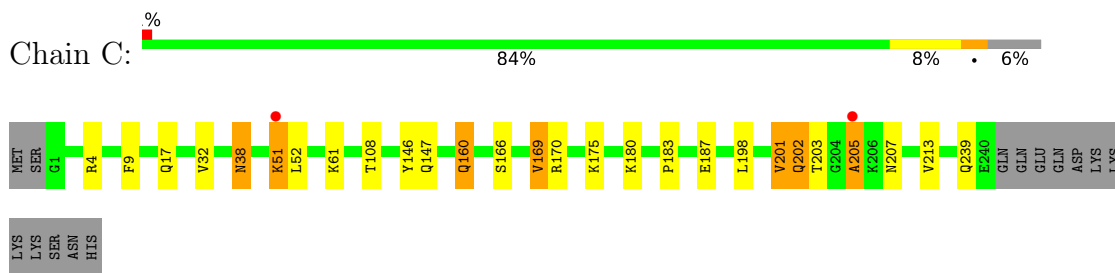
- Molecule 2: Proteasome subunit alpha type-3



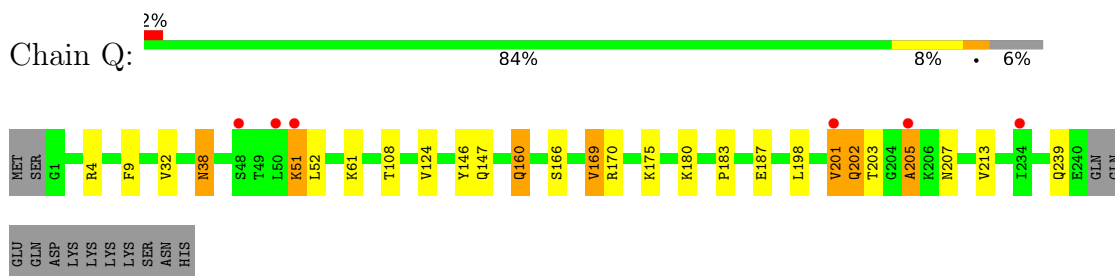
- Molecule 2: Proteasome subunit alpha type-3



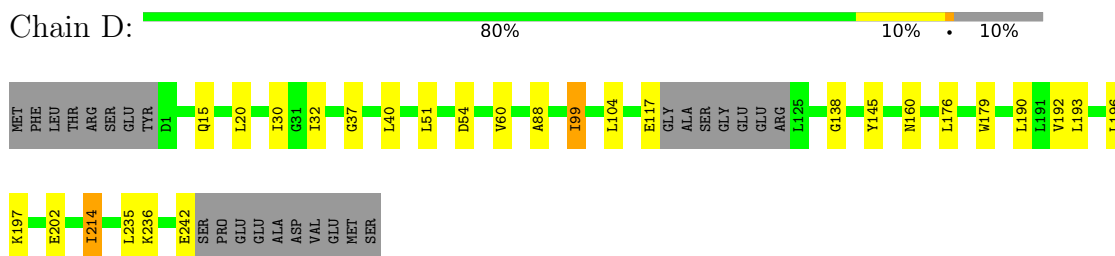
- Molecule 3: Proteasome subunit alpha type-4



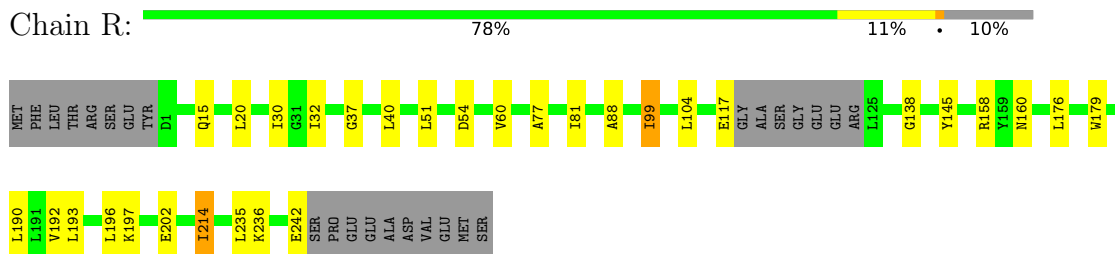
- Molecule 3: Proteasome subunit alpha type-4



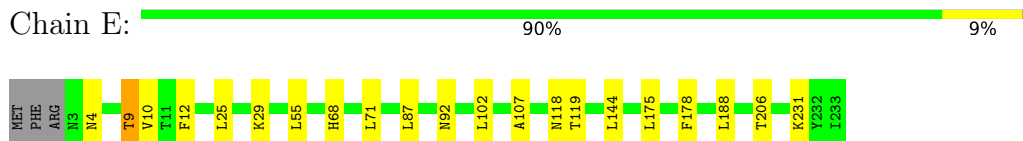
- Molecule 4: Proteasome subunit alpha type-5



- Molecule 4: Proteasome subunit alpha type-5



- Molecule 5: Proteasome subunit alpha type-6



- Molecule 5: Proteasome subunit alpha type-6



- Molecule 9: Proteasome subunit beta type-3

Chain I:  90% 9%



- Molecule 9: Proteasome subunit beta type-3

Chain W:  90% 9%



- Molecule 10: Proteasome subunit beta type-4

Chain J:  87% 10% ..



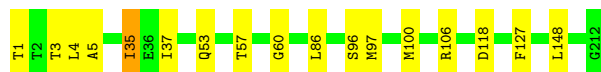
- Molecule 10: Proteasome subunit beta type-4

Chain X:  87% 10% ..



- Molecule 11: Proteasome subunit beta type-5

Chain K:  92% 8%



- Molecule 11: Proteasome subunit beta type-5

Chain Y:  91% 8%



- Molecule 12: Proteasome subunit beta type-6

Chain L:  91% 8%



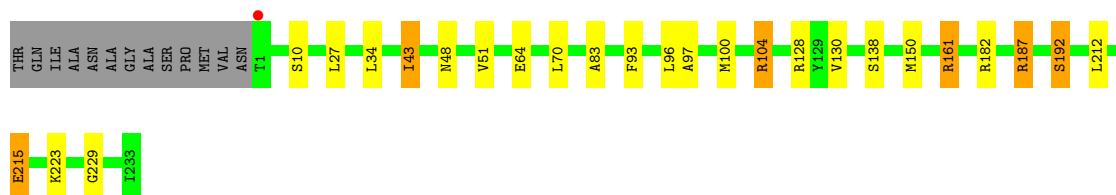
- Molecule 12: Proteasome subunit beta type-6

Chain Z:  89% 10%



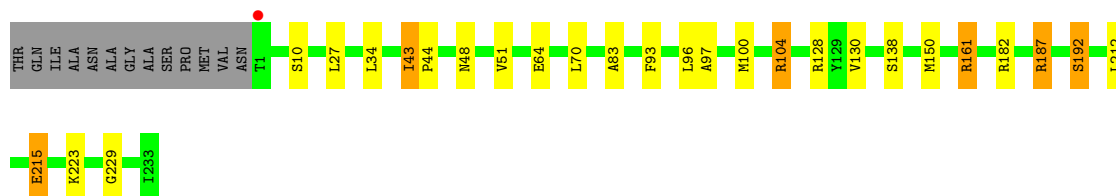
- Molecule 13: Proteasome subunit beta type-7

Chain M:  84% 8% 5%

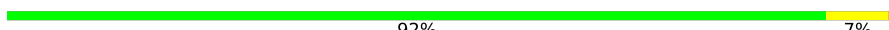


- Molecule 13: Proteasome subunit beta type-7

Chain a:  84% 9% 5%



- Molecule 14: Proteasome subunit beta type-1

Chain N:  92% 7%



- Molecule 14: Proteasome subunit beta type-1

Chain b:  92% 7%



- Molecule 15: Ac-LAE-ep

Chain c:  60% 40%

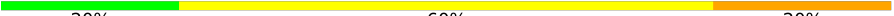


- Molecule 15: Ac-LAE-ep

Chain d:  80% 20%



• Molecule 15: Ac-LAE-ep

Chain e:  20% 60% 20%



• Molecule 15: Ac-LAE-ep

Chain f:  80% 20%



• Molecule 15: Ac-LAE-ep

Chain g:  60% 20% 20%



• Molecule 15: Ac-LAE-ep

Chain h:  60% 40%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	136.35Å 300.08Å 144.78Å 90.00° 112.99° 90.00°	Depositor
Resolution (Å)	15.00 – 2.70 15.00 – 2.70	Depositor EDS
% Data completeness (in resolution range)	97.2 (15.00-2.70) 96.7 (15.00-2.70)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.63 (at 2.69Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.188 , 0.221 0.193 , 0.225	Depositor DCC
R_{free} test set	14123 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	63.6	Xtriage
Anisotropy	0.035	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 57.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	49825	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.48% 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: POL, ACE, GAU, MG, MES

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	0.39	0/1952	0.71	0/2642
1	O	0.39	0/1952	0.71	0/2642
2	B	0.40	0/1934	0.71	0/2618
2	P	0.39	0/1934	0.71	0/2618
3	C	0.39	0/1910	0.74	0/2586
3	Q	0.39	0/1910	0.74	0/2586
4	D	0.39	0/1837	0.70	0/2475
4	R	0.39	0/1837	0.70	0/2475
5	E	0.39	0/1800	0.68	0/2433
5	S	0.39	0/1800	0.68	0/2433
6	F	0.40	0/1932	0.70	0/2609
6	T	0.40	0/1932	0.71	0/2609
7	G	0.37	0/1945	0.70	0/2634
7	U	0.38	0/1945	0.70	0/2634
8	H	0.44	0/1712	0.70	0/2322
8	V	0.39	0/1723	0.79	5/2337 (0.2%)
9	I	0.36	0/1611	0.71	0/2174
9	W	0.36	0/1611	0.71	0/2174
10	J	0.39	0/1589	0.69	0/2142
10	X	0.38	0/1589	0.69	0/2142
11	K	0.41	0/1681	0.70	0/2274
11	Y	0.40	0/1681	0.69	0/2274
12	L	0.37	0/1795	0.67	0/2420
12	Z	0.38	0/1806	0.67	0/2435
13	M	0.39	0/1855	0.69	0/2514
13	a	0.39	0/1855	0.70	0/2514
14	N	0.48	0/1541	0.72	0/2087
14	b	0.45	0/1541	0.72	0/2087
15	c	1.23	0/13	0.98	0/17
15	d	1.18	0/13	1.15	0/17
15	e	1.10	0/13	0.94	0/17
15	f	0.66	0/13	0.99	0/17

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
15	g	0.81	0/13	1.02	0/17
15	h	1.27	0/13	1.18	0/17
All	All	0.40	0/50288	0.71	5/67992 (0.0%)

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
8	V	0	1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	V	116	ASP	N-CA-C	-9.88	102.02	114.56
8	V	115	ALA	CA-C-N	9.53	136.66	122.08
8	V	115	ALA	C-N-CA	9.53	136.66	122.08
8	V	116	ASP	N-CA-CB	-5.84	103.23	111.00
8	V	115	ALA	N-CA-C	5.62	119.75	113.01

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
8	V	115	ALA	Peptide

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	4	0
1	O	1915	0	1929	5	0
2	B	1904	0	1904	17	0
2	P	1904	0	1904	16	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	1881	0	1895	13	0
3	Q	1881	0	1895	14	0
4	D	1813	0	1797	8	0
4	R	1813	0	1797	10	0
5	E	1773	0	1775	5	0
5	S	1773	0	1775	6	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	8	0
8	H	1682	0	1682	7	0
8	V	1689	0	1689	6	0
9	I	1581	0	1574	11	0
9	W	1581	0	1574	9	0
10	J	1561	0	1569	12	0
10	X	1561	0	1569	11	0
11	K	1644	0	1592	9	0
11	Y	1644	0	1592	11	0
12	L	1757	0	1711	11	0
12	Z	1764	0	1718	15	0
13	M	1824	0	1832	11	0
13	a	1824	0	1832	12	0
14	N	1512	0	1478	10	0
14	b	1512	0	1478	10	0
15	c	29	0	30	1	0
15	d	29	0	30	0	0
15	e	29	0	30	4	0
15	f	29	0	30	0	0
15	g	29	0	30	1	0
15	h	29	0	30	1	0
16	G	1	0	0	0	0
16	H	1	0	0	0	0
16	I	2	0	0	0	0
16	K	1	0	0	0	0
16	L	1	0	0	0	0
16	N	1	0	0	0	0
17	K	12	0	13	0	0
17	Y	12	0	13	0	0
18	A	6	0	0	0	0
18	B	13	0	0	0	0
18	C	10	0	0	0	0
18	D	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
18	E	5	0	0	0	0
18	F	13	0	0	0	0
18	G	15	0	0	0	0
18	H	23	0	0	0	0
18	I	10	0	0	0	0
18	J	11	0	0	0	0
18	K	14	0	0	0	0
18	L	14	0	0	0	0
18	M	13	0	0	0	0
18	N	6	0	0	0	0
18	O	3	0	0	0	0
18	P	7	0	0	0	0
18	Q	8	0	0	0	0
18	R	3	0	0	0	0
18	S	4	0	0	0	0
18	T	9	0	0	0	0
18	U	21	0	0	0	0
18	V	13	0	0	0	0
18	W	8	0	0	0	0
18	X	14	0	0	0	0
18	Y	16	0	0	0	0
18	Z	20	0	0	0	0
18	a	17	0	0	0	0
18	b	11	0	0	0	0
18	c	1	0	0	0	0
18	d	1	0	0	0	0
18	e	1	0	0	0	0
18	f	1	0	0	0	0
18	g	1	0	0	0	0
All	All	49825	0	49264	242	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 (242) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:K:100:MET:HE3	11:K:127:PHE:HB2	1.56	0.88
11:Y:100:MET:HE3	11:Y:127:PHE:HB2	1.55	0.87
4:R:99:ILE:HD11	4:R:104:LEU:HB2	1.57	0.86
4:D:99:ILE:HD11	4:D:104:LEU:HB2	1.57	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:V:116:ASP:OD2	14:b:50:ALA:HB2	1.80	0.81
12:Z:108[B]:HIS:HD2	12:Z:139:GLY:HA3	1.48	0.79
11:K:53:GLN:O	11:K:57:THR:HG23	1.86	0.75
11:Y:53:GLN:O	11:Y:57:THR:HG23	1.85	0.74
12:Z:108[B]:HIS:CD2	12:Z:139:GLY:HA3	2.23	0.74
8:V:116:ASP:OD2	14:b:50:ALA:CB	2.37	0.72
10:J:16:ALA:HB2	10:J:161:LEU:HD21	1.78	0.66
10:X:16:ALA:HB2	10:X:161:LEU:HD21	1.78	0.65
8:H:168:GLY:O	15:c:5:POL:H33	1.99	0.62
4:D:88:ALA:HA	4:D:99:ILE:HG21	1.82	0.62
3:C:160:GLN:HE21	3:C:160:GLN:HA	1.65	0.61
3:Q:160:GLN:HA	3:Q:160:GLN:HE21	1.65	0.61
8:V:116:ASP:OD2	8:V:116:ASP:C	2.44	0.61
4:R:88:ALA:HA	4:R:99:ILE:HG21	1.82	0.61
14:b:152:VAL:HA	14:b:175:MET:HE1	1.82	0.61
14:N:152:VAL:HA	14:N:175:MET:HE1	1.83	0.61
14:b:22:THR:HG23	14:b:22:THR:O	2.00	0.61
7:U:23:PHE:O	7:U:26:THR:HB	2.03	0.59
14:N:13:ILE:HG21	14:N:175:MET:HE2	1.85	0.58
14:N:136:GLY:HA2	14:b:161:GLN:HE21	1.67	0.58
7:G:23:PHE:O	7:G:26:THR:HB	2.03	0.58
14:N:161:GLN:HE21	14:b:136:GLY:HA2	1.67	0.58
14:b:13:ILE:HG21	14:b:175:MET:HE2	1.86	0.58
3:C:51:LYS:O	3:C:52:LEU:HB2	2.06	0.56
12:L:23:LEU:HD13	12:L:43:VAL:HG13	1.88	0.56
2:B:93:HIS:HB3	2:B:113:ARG:HH21	1.70	0.56
12:Z:13:LEU:HD13	12:Z:150:LEU:HD21	1.89	0.55
3:Q:51:LYS:O	3:Q:52:LEU:HB2	2.06	0.55
9:W:10:ILE:HG21	9:W:141:ALA:HB3	1.89	0.55
2:P:93:HIS:HB3	2:P:113:ARG:HH21	1.71	0.55
12:L:13:LEU:HD13	12:L:150:LEU:HD21	1.88	0.54
15:e:3:ALA:C	15:e:4:GAU:HG2	2.31	0.54
10:X:1:MET:HA	10:X:34:LYS:HE3	1.90	0.54
11:Y:5:ALA:HB3	11:Y:100:MET:HE2	1.89	0.54
8:H:52:THR:O	8:H:56:THR:HB	2.08	0.54
9:I:10:ILE:HG21	9:I:141:ALA:HB3	1.89	0.54
5:S:87:LEU:HD21	5:S:107:ALA:HB1	1.91	0.53
2:P:183:MET:HE1	2:P:191:LEU:HD12	1.91	0.53
12:Z:23:LEU:HD13	12:Z:43:VAL:HG13	1.88	0.53
10:J:1:MET:HA	10:J:34:LYS:HE3	1.90	0.53
11:K:5:ALA:HB3	11:K:100:MET:HE2	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:V:52:THR:O	8:V:56:THR:HB	2.09	0.53
11:Y:35:ILE:HD13	11:Y:37:ILE:HG13	1.90	0.52
12:L:8:ASN:HA	12:L:30:ILE:O	2.09	0.52
2:B:183:MET:HE1	2:B:191:LEU:HD12	1.90	0.52
9:W:20:VAL:HG13	9:W:118:PRO:HB3	1.90	0.52
5:E:87:LEU:HD21	5:E:107:ALA:HB1	1.91	0.52
11:K:35:ILE:HD13	11:K:37:ILE:HG13	1.90	0.51
14:N:22:THR:HG23	14:N:22:THR:O	2.09	0.51
12:Z:8:ASN:HA	12:Z:30:ILE:O	2.10	0.51
4:R:138:GLY:HA2	4:R:214:ILE:HG12	1.93	0.51
9:I:20:VAL:HG13	9:I:118:PRO:HB3	1.91	0.51
13:M:43:ILE:HG21	13:M:64:GLU:HG2	1.93	0.51
11:K:5:ALA:HB3	11:K:100:MET:CE	2.41	0.51
11:Y:5:ALA:HB3	11:Y:100:MET:CE	2.41	0.51
8:H:112:SER:OG	8:H:120:ASP:HB2	2.12	0.50
9:W:9:GLY:HA3	9:W:41:LYS:HE2	1.93	0.50
10:J:174:MET:HA	10:X:174:MET:HA	1.92	0.50
7:G:34:LEU:C	7:G:34:LEU:HD23	2.37	0.50
10:X:1:MET:HA	10:X:34:LYS:CE	2.41	0.50
11:Y:1:THR:CG2	11:Y:3:THR:HG23	2.42	0.50
13:a:43:ILE:HG21	13:a:64:GLU:HG2	1.93	0.50
10:J:1:MET:HA	10:J:34:LYS:CE	2.42	0.49
4:D:138:GLY:HA2	4:D:214:ILE:HG12	1.93	0.49
9:I:9:GLY:HA3	9:I:41:LYS:HE2	1.93	0.49
7:U:34:LEU:C	7:U:34:LEU:HD23	2.37	0.49
9:I:65:MET:HE1	9:I:93:SER:HB3	1.95	0.49
12:L:13:LEU:CD1	12:L:150:LEU:HD21	2.42	0.49
5:E:12:PHE:H	6:F:19:GLN:HE22	1.58	0.49
9:I:101:PRO:HB3	9:I:126:ILE:HD12	1.94	0.49
9:W:65:MET:HE1	9:W:93:SER:HB3	1.95	0.49
11:K:37:ILE:HG23	11:K:60:GLY:HA2	1.95	0.49
1:O:12:PHE:H	2:P:20:GLN:HE22	1.61	0.49
11:Y:37:ILE:HG23	11:Y:60:GLY:HA2	1.95	0.48
5:S:12:PHE:H	6:T:19:GLN:HE22	1.60	0.48
9:W:101:PRO:HB3	9:W:126:ILE:HD12	1.95	0.48
12:L:31:THR:HG23	12:L:36:ASN:HD21	1.78	0.48
10:X:152:MET:HE2	10:X:156:GLU:HB3	1.94	0.48
12:Z:31:THR:HG23	12:Z:36:ASN:HD21	1.78	0.48
12:L:4:PRO:O	13:M:104:ARG:NH1	2.43	0.48
3:Q:201:VAL:O	3:Q:202:GLN:CB	2.62	0.48
6:T:123:ASN:C	6:T:123:ASN:HD22	2.22	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:Z:13:LEU:CD1	12:Z:150:LEU:HD21	2.43	0.48
10:J:152:MET:HE2	10:J:156:GLU:HB3	1.95	0.47
3:C:201:VAL:O	3:C:202:GLN:CB	2.62	0.47
5:S:68:HIS:HE1	5:S:102:LEU:O	1.97	0.47
6:T:155:GLY:HA3	7:U:59:THR:HG21	1.97	0.47
14:N:22:THR:HB	15:e:2:LEU:HD23	1.96	0.47
11:Y:25:TRP:CH2	12:Z:144:SER:HA	2.50	0.47
2:B:180:LYS:O	2:B:183:MET:HB2	2.14	0.47
2:P:180:LYS:O	2:P:183:MET:HB2	2.14	0.47
5:E:68:HIS:HE1	5:E:102:LEU:O	1.96	0.47
4:D:32:ILE:HD12	4:D:192:VAL:HG23	1.97	0.47
3:Q:9:PHE:H	4:R:15:GLN:HE22	1.63	0.47
7:U:83:ASN:C	7:U:83:ASN:HD22	2.23	0.47
1:O:122:THR:HG22	2:P:128:ARG:HH21	1.79	0.47
4:R:32:ILE:HD12	4:R:192:VAL:HG23	1.96	0.47
10:X:3:ILE:HD12	10:X:176:PHE:CG	2.50	0.47
12:Z:195:HIS:HD2	12:Z:197:GLN:H	1.63	0.47
3:C:201:VAL:O	3:C:202:GLN:HB2	2.14	0.46
3:C:201:VAL:HG13	3:C:202:GLN:H	1.80	0.46
6:F:123:ASN:C	6:F:123:ASN:HD22	2.23	0.46
10:J:126:VAL:HG12	10:J:128:LEU:HG	1.98	0.46
1:O:23:TYR:CD1	7:U:12:PRO:HA	2.50	0.46
4:R:160:ASN:HB3	4:R:179:TRP:CE2	2.51	0.46
4:D:37:GLY:HA2	4:D:145:TYR:CE1	2.51	0.46
10:X:126:VAL:HG12	10:X:128:LEU:HG	1.98	0.46
3:Q:201:VAL:HG13	3:Q:202:GLN:H	1.80	0.46
9:W:101:PRO:HB3	9:W:126:ILE:CD1	2.46	0.46
3:Q:201:VAL:O	3:Q:202:GLN:HB2	2.14	0.46
1:A:115:ALA:HB1	1:A:154:GLY:O	2.16	0.46
12:L:195:HIS:HD2	12:L:197:GLN:H	1.63	0.46
8:H:80:LEU:HD12	8:H:113:ILE:HD11	1.97	0.46
4:D:160:ASN:HB3	4:D:179:TRP:CE2	2.51	0.45
7:G:83:ASN:C	7:G:83:ASN:HD22	2.24	0.45
9:I:36:SER:HB2	10:J:126:VAL:HG11	1.96	0.45
4:R:37:GLY:HA2	4:R:145:TYR:CE1	2.51	0.45
6:F:172:LEU:HD13	6:F:195:ILE:HD13	1.99	0.45
1:O:115:ALA:HB1	1:O:154:GLY:O	2.17	0.45
10:J:3:ILE:HD12	10:J:176:PHE:CG	2.50	0.45
13:M:161:ARG:HG3	13:M:161:ARG:HH11	1.82	0.45
7:G:78:ILE:N	7:G:79:PRO:CD	2.80	0.45
13:M:182:ARG:NH2	13:M:215:GLU:O	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:T:172:LEU:HD13	6:T:195:ILE:HD13	1.99	0.45
13:a:161:ARG:HG3	13:a:161:ARG:HH11	1.82	0.45
9:I:101:PRO:HB3	9:I:126:ILE:CD1	2.46	0.45
6:F:172:LEU:CD1	6:F:195:ILE:HD13	2.47	0.45
8:H:112:SER:HG	8:H:120:ASP:HB2	1.82	0.45
5:S:175:LEU:HA	5:S:178:PHE:CE2	2.52	0.45
5:E:9:THR:HG21	5:E:119:THR:HA	1.99	0.45
10:X:3:ILE:HD12	10:X:176:PHE:CD2	2.52	0.45
12:Z:4:PRO:O	13:a:104:ARG:NH1	2.44	0.45
5:E:175:LEU:HA	5:E:178:PHE:CE2	2.52	0.44
3:C:205:ALA:C	3:C:207:ASN:H	2.26	0.44
12:L:17:GLY:HA3	12:L:20:PHE:CE1	2.53	0.44
4:R:30:ILE:HD12	4:R:196:LEU:HG	2.00	0.44
6:T:172:LEU:CD1	6:T:195:ILE:HD13	2.47	0.44
12:Z:17:GLY:HA3	12:Z:20:PHE:CE1	2.53	0.44
4:D:30:ILE:HD12	4:D:196:LEU:HG	2.00	0.44
9:W:36:SER:HB2	10:X:126:VAL:HG11	2.00	0.44
1:A:122:THR:CG2	2:B:128:ARG:HH21	2.31	0.44
1:O:122:THR:CG2	2:P:128:ARG:HH21	2.30	0.44
8:H:104:ASP:HB2	8:H:105:PRO:HD2	1.99	0.44
10:J:3:ILE:HD12	10:J:176:PHE:CD2	2.52	0.44
14:N:38:HIS:O	14:N:39:ASP:C	2.60	0.44
2:P:47:ALA:HB1	2:P:64:LYS:HD2	2.00	0.44
3:Q:205:ALA:C	3:Q:207:ASN:H	2.26	0.44
13:a:96:LEU:O	13:a:100:MET:HG2	2.18	0.44
14:b:38:HIS:O	14:b:39:ASP:C	2.60	0.44
11:K:1:THR:CG2	11:K:3:THR:HG23	2.48	0.44
13:M:150:MET:HE1	13:M:187:ARG:HG2	2.00	0.44
2:B:47:ALA:HB1	2:B:64:LYS:HD2	2.00	0.43
13:M:43:ILE:HG12	13:M:51:VAL:HB	2.00	0.43
5:S:9:THR:HG21	5:S:119:THR:HA	1.99	0.43
12:Z:49:ASN:HD21	12:Z:211:GLY:HA2	1.83	0.43
13:a:27:LEU:HD21	13:a:34:LEU:HD22	2.00	0.43
13:a:43:ILE:HG12	13:a:51:VAL:HB	2.00	0.43
2:P:151:ASN:HB2	2:P:152:PRO:CD	2.48	0.43
7:U:83:ASN:C	7:U:83:ASN:ND2	2.76	0.43
8:V:104:ASP:HB2	8:V:105:PRO:HD2	1.99	0.43
2:B:50:LYS:O	2:B:51:VAL:C	2.62	0.43
12:L:49:ASN:HD21	12:L:211:GLY:HA2	1.83	0.43
13:M:96:LEU:O	13:M:100:MET:HG2	2.18	0.43
7:U:78:ILE:N	7:U:79:PRO:CD	2.81	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:Q:160:GLN:HE22	3:Q:170:ARG:HE	1.67	0.43
14:b:176:VAL:HG12	14:b:178:LEU:HD13	2.01	0.43
2:P:50:LYS:O	2:P:51:VAL:C	2.61	0.43
13:a:182:ARG:NH2	13:a:215:GLU:O	2.51	0.43
1:A:12:PHE:H	2:B:20:GLN:HE22	1.67	0.43
2:B:151:ASN:HB2	2:B:152:PRO:CD	2.48	0.43
3:C:160:GLN:HE22	3:C:170:ARG:HE	1.67	0.43
2:B:12:PHE:H	3:C:17:GLN:HE22	1.67	0.43
13:a:27:LEU:HB2	13:a:192:SER:HB3	2.01	0.43
13:a:150:MET:HE1	13:a:187:ARG:HG2	2.00	0.43
12:L:62:ASP:O	12:L:66:LYS:HB2	2.19	0.42
13:M:93:PHE:CE2	13:M:128:ARG:HD3	2.54	0.42
10:J:3:ILE:HB	10:J:18:SER:HB3	2.01	0.42
13:M:27:LEU:HD21	13:M:34:LEU:HD22	2.01	0.42
2:P:151:ASN:HB2	2:P:152:PRO:HD2	2.01	0.42
13:a:93:PHE:CE2	13:a:128:ARG:HD3	2.55	0.42
1:A:122:THR:HG22	2:B:128:ARG:HH21	1.84	0.42
11:Y:86:LEU:C	11:Y:86:LEU:HD13	2.45	0.42
14:b:22:THR:HB	15:h:2:LEU:CD2	2.49	0.42
3:C:9:PHE:H	4:D:15:GLN:HE22	1.67	0.42
7:G:83:ASN:C	7:G:83:ASN:ND2	2.77	0.42
3:C:108:THR:HG21	3:C:146:TYR:HB3	2.02	0.42
10:J:36:ARG:NH1	10:J:58:GLU:OE2	2.52	0.42
2:P:36:GLY:C	2:P:147:LEU:HD11	2.44	0.42
2:P:124:HIS:HB3	3:Q:124:VAL:HG12	2.01	0.42
12:Z:62:ASP:O	12:Z:66:LYS:HB2	2.19	0.42
2:B:110:LEU:HD23	2:B:110:LEU:C	2.45	0.42
2:B:204:ALA:O	2:B:209:ARG:NH2	2.52	0.42
2:B:151:ASN:HB2	2:B:152:PRO:HD2	2.02	0.42
2:P:204:ALA:O	2:P:209:ARG:NH2	2.53	0.42
4:R:77:ALA:O	4:R:81:ILE:HG12	2.20	0.42
3:C:38:ASN:N	3:C:38:ASN:HD22	2.18	0.42
9:I:14:MET:HB3	9:I:162:LEU:HD11	2.02	0.42
11:K:86:LEU:C	11:K:86:LEU:HD13	2.45	0.41
14:N:176:VAL:HG12	14:N:178:LEU:HD13	2.01	0.41
8:V:206:PRO:O	8:V:209:THR:OG1	2.31	0.41
2:B:36:GLY:C	2:B:147:LEU:HD11	2.45	0.41
3:C:166:SER:HA	3:C:169:VAL:HG13	2.02	0.41
14:N:22:THR:HB	15:e:2:LEU:CD2	2.50	0.41
2:P:57:GLU:O	2:P:61:SER:HB2	2.20	0.41
3:Q:38:ASN:N	3:Q:38:ASN:HD22	2.18	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:M:27:LEU:HB2	13:M:192:SER:HB3	2.02	0.41
10:X:36:ARG:NH1	10:X:58:GLU:OE2	2.54	0.41
9:I:98:ARG:O	9:I:126:ILE:HD11	2.21	0.41
2:P:110:LEU:C	2:P:110:LEU:HD23	2.45	0.41
7:U:63:ILE:HD12	7:U:215:GLU:HG2	2.02	0.41
9:W:14:MET:HB3	9:W:162:LEU:HD11	2.03	0.41
13:a:97:ALA:HA	13:a:130:VAL:HG21	2.03	0.41
2:B:57:GLU:O	2:B:61:SER:HB2	2.21	0.41
13:M:97:ALA:HA	13:M:130:VAL:HG21	2.03	0.41
3:Q:166:SER:HA	3:Q:169:VAL:HG13	2.02	0.41
9:I:26:LEU:HD21	9:I:185:VAL:HG23	2.02	0.41
12:L:136:CYS:SG	12:L:150:LEU:HB3	2.61	0.41
3:Q:198:LEU:HA	3:Q:201:VAL:HG12	2.02	0.41
2:P:139:TYR:CD1	2:P:224:VAL:HG21	2.56	0.41
3:Q:108:THR:HG21	3:Q:146:TYR:HB3	2.02	0.41
3:Q:201:VAL:HG13	3:Q:202:GLN:N	2.36	0.41
11:Y:52:CYS:HA	11:Y:97:MET:HE3	2.03	0.41
10:J:67:TYR:CE1	10:J:75:LEU:HD13	2.56	0.41
8:H:50:ALA:CB	9:I:126:ILE:HG23	2.52	0.40
11:K:96:SER:C	11:K:97:MET:HG3	2.46	0.40
2:B:139:TYR:CD1	2:B:224:VAL:HG21	2.56	0.40
4:R:158:ARG:HB3	5:S:57:SER:HB3	2.03	0.40
10:X:67:TYR:CE1	10:X:75:LEU:HD13	2.56	0.40
11:Y:5:ALA:HA	11:Y:13:ILE:O	2.22	0.40
12:Z:126:ASP:OD1	12:Z:126:ASP:C	2.65	0.40
13:a:43:ILE:HA	13:a:44:PRO:HD3	1.94	0.40
2:B:46:ALA:HB2	2:B:212:PHE:CE1	2.57	0.40
14:N:168:SER:HB3	15:e:5:POL:H33	2.03	0.40
12:Z:18:GLU:O	12:Z:119:LYS:HA	2.22	0.40
7:G:63:ILE:HD12	7:G:215:GLU:HG2	2.03	0.40
9:W:26:LEU:HD21	9:W:185:VAL:HG23	2.02	0.40
3:C:198:LEU:HA	3:C:201:VAL:HG12	2.02	0.40
15:g:4:GAU:N	15:g:5:POL:H12	2.36	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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%)	237 (96%)	11 (4%)	0	100	100
1	O	248/250 (99%)	237 (96%)	11 (4%)	0	100	100
2	B	242/258 (94%)	231 (96%)	9 (4%)	2 (1%)	16	37
2	P	242/258 (94%)	231 (96%)	9 (4%)	2 (1%)	16	37
3	C	238/254 (94%)	229 (96%)	4 (2%)	5 (2%)	5	15
3	Q	238/254 (94%)	229 (96%)	4 (2%)	5 (2%)	5	15
4	D	231/260 (89%)	225 (97%)	6 (3%)	0	100	100
4	R	231/260 (89%)	225 (97%)	6 (3%)	0	100	100
5	E	229/234 (98%)	218 (95%)	11 (5%)	0	100	100
5	S	229/234 (98%)	218 (95%)	11 (5%)	0	100	100
6	F	241/288 (84%)	234 (97%)	7 (3%)	0	100	100
6	T	241/288 (84%)	234 (97%)	7 (3%)	0	100	100
7	G	239/252 (95%)	234 (98%)	5 (2%)	0	100	100
7	U	239/252 (95%)	234 (98%)	5 (2%)	0	100	100
8	H	220/232 (95%)	214 (97%)	6 (3%)	0	100	100
8	V	221/232 (95%)	216 (98%)	5 (2%)	0	100	100
9	I	202/205 (98%)	196 (97%)	6 (3%)	0	100	100
9	W	202/205 (98%)	196 (97%)	6 (3%)	0	100	100
10	J	193/198 (98%)	188 (97%)	4 (2%)	1 (0%)	24	48
10	X	193/198 (98%)	188 (97%)	4 (2%)	1 (0%)	24	48
11	K	210/212 (99%)	203 (97%)	7 (3%)	0	100	100
11	Y	210/212 (99%)	203 (97%)	7 (3%)	0	100	100
12	L	220/222 (99%)	216 (98%)	4 (2%)	0	100	100
12	Z	221/222 (100%)	217 (98%)	4 (2%)	0	100	100
13	M	231/246 (94%)	223 (96%)	6 (3%)	2 (1%)	14	35

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
13	a	231/246 (94%)	223 (96%)	6 (3%)	2 (1%)	14	35
14	N	194/196 (99%)	186 (96%)	8 (4%)	0	100	100
14	b	194/196 (99%)	186 (96%)	8 (4%)	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
15	h	2/5 (40%)	2 (100%)	0	0	100	100
All	All	6290/6644 (95%)	6083 (97%)	187 (3%)	20 (0%)	36	60

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	202	GLN
3	Q	202	GLN
2	B	51	VAL
3	C	239	GLN
10	J	2	ASP
2	P	51	VAL
3	Q	239	GLN
10	X	2	ASP
3	C	183	PRO
3	Q	183	PRO
2	B	221	ASP
3	C	205	ALA
2	P	221	ASP
3	Q	205	ALA
13	M	83	ALA
13	a	83	ALA
3	C	201	VAL
3	Q	201	VAL
13	M	229	GLY
13	a	229	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	209/209 (100%)	203 (97%)	6 (3%)	37	67
1	O	209/209 (100%)	203 (97%)	6 (3%)	37	67
2	B	203/216 (94%)	193 (95%)	10 (5%)	22	49
2	P	203/216 (94%)	193 (95%)	10 (5%)	22	49
3	C	212/226 (94%)	199 (94%)	13 (6%)	17	40
3	Q	212/226 (94%)	199 (94%)	13 (6%)	17	40
4	D	194/215 (90%)	178 (92%)	16 (8%)	10	27
4	R	194/215 (90%)	178 (92%)	16 (8%)	10	27
5	E	190/193 (98%)	177 (93%)	13 (7%)	14	35
5	S	190/193 (98%)	177 (93%)	13 (7%)	14	35
6	F	201/239 (84%)	190 (94%)	11 (6%)	19	45
6	T	201/239 (84%)	190 (94%)	11 (6%)	19	45
7	G	206/210 (98%)	195 (95%)	11 (5%)	20	46
7	U	206/210 (98%)	195 (95%)	11 (5%)	20	46
8	H	181/190 (95%)	173 (96%)	8 (4%)	25	53
8	V	182/190 (96%)	174 (96%)	8 (4%)	25	53
9	I	172/173 (99%)	167 (97%)	5 (3%)	37	67
9	W	172/173 (99%)	167 (97%)	5 (3%)	37	67
10	J	173/175 (99%)	166 (96%)	7 (4%)	28	56
10	X	173/175 (99%)	165 (95%)	8 (5%)	24	51
11	K	169/169 (100%)	164 (97%)	5 (3%)	36	66
11	Y	169/169 (100%)	164 (97%)	5 (3%)	36	66
12	L	185/185 (100%)	180 (97%)	5 (3%)	39	69
12	Z	186/185 (100%)	181 (97%)	5 (3%)	39	69
13	M	199/208 (96%)	187 (94%)	12 (6%)	17	41
13	a	199/208 (96%)	187 (94%)	12 (6%)	17	41

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
14	N	162/162 (100%)	157 (97%)	5 (3%)	35	65
14	b	162/162 (100%)	157 (97%)	5 (3%)	35	65
15	c	1/1 (100%)	1 (100%)	0	100	100
15	d	1/1 (100%)	1 (100%)	0	100	100
15	e	1/1 (100%)	1 (100%)	0	100	100
15	f	1/1 (100%)	1 (100%)	0	100	100
15	g	1/1 (100%)	1 (100%)	0	100	100
15	h	1/1 (100%)	1 (100%)	0	100	100
All	All	5320/5546 (96%)	5065 (95%)	255 (5%)	23	50

All (255) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	LYS
1	A	59	GLU
1	A	61	LEU
1	A	122	THR
1	A	157	PHE
1	A	250	LEU
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	180	LYS
2	B	184	LYS
2	B	191	LEU
3	C	4	ARG
3	C	32	VAL
3	C	38	ASN
3	C	51	LYS
3	C	61	LYS
3	C	147	GLN
3	C	160	GLN
3	C	169	VAL
3	C	175	LYS
3	C	180	LYS

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Mol	Chain	Res	Type
3	C	187	GLU
3	C	203	THR
3	C	213	VAL
4	D	20	LEU
4	D	40	LEU
4	D	51	LEU
4	D	54	ASP
4	D	60	VAL
4	D	99	ILE
4	D	117	GLU
4	D	176	LEU
4	D	190	LEU
4	D	193	LEU
4	D	197	LYS
4	D	202	GLU
4	D	214	ILE
4	D	235	LEU
4	D	236	LYS
4	D	242	GLU
5	E	4	ASN
5	E	9	THR
5	E	10	VAL
5	E	25	LEU
5	E	29	LYS
5	E	55	LEU
5	E	71	LEU
5	E	92	ASN
5	E	118	ASN
5	E	144	LEU
5	E	188	LEU
5	E	206	THR
5	E	231	LYS
6	F	14	ASP
6	F	68	ARG
6	F	94	SER
6	F	123	ASN
6	F	165	ARG
6	F	172	LEU
6	F	181	GLU
6	F	187	GLU
6	F	203	ASN
6	F	206	LYS

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Mol	Chain	Res	Type
6	F	221	ASN
7	G	26	THR
7	G	28	GLN
7	G	34	LEU
7	G	83	ASN
7	G	115	LEU
7	G	125	MET
7	G	166	GLN
7	G	171	THR
7	G	181	LYS
7	G	235	ARG
7	G	236	LEU
8	H	7	LYS
8	H	13	VAL
8	H	34	LEU
8	H	56	THR
8	H	68	LEU
8	H	127	LEU
8	H	196	ARG
8	H	198	GLU
9	I	37	ASN
9	I	126	ILE
9	I	171	LEU
9	I	182	TRP
9	I	192	ASP
10	J	2	ASP
10	J	23	ARG
10	J	35	THR
10	J	75	LEU
10	J	91	SER
10	J	144	LEU
10	J	163	LEU
11	K	4	LEU
11	K	35	ILE
11	K	106	ARG
11	K	118	ASP
11	K	148	LEU
12	L	1	GLN
12	L	23	LEU
12	L	49	ASN
12	L	150	LEU
12	L	167	LYS

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Mol	Chain	Res	Type
13	M	10	SER
13	M	43	ILE
13	M	48	ASN
13	M	70	LEU
13	M	104	ARG
13	M	138	SER
13	M	161	ARG
13	M	187	ARG
13	M	192	SER
13	M	212	LEU
13	M	215	GLU
13	M	223	LYS
14	N	9	LYS
14	N	21	THR
14	N	36	ARG
14	N	119	VAL
14	N	178	LEU
1	O	17	LYS
1	O	59	GLU
1	O	61	LEU
1	O	122	THR
1	O	157	PHE
1	O	250	LEU
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	102	ASN
2	P	180	LYS
2	P	184	LYS
2	P	191	LEU
3	Q	4	ARG
3	Q	32	VAL
3	Q	38	ASN
3	Q	51	LYS
3	Q	61	LYS
3	Q	147	GLN
3	Q	160	GLN
3	Q	169	VAL
3	Q	175	LYS

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Mol	Chain	Res	Type
3	Q	180	LYS
3	Q	187	GLU
3	Q	203	THR
3	Q	213	VAL
4	R	20	LEU
4	R	40	LEU
4	R	51	LEU
4	R	54	ASP
4	R	60	VAL
4	R	99	ILE
4	R	117	GLU
4	R	176	LEU
4	R	190	LEU
4	R	193	LEU
4	R	197	LYS
4	R	202	GLU
4	R	214	ILE
4	R	235	LEU
4	R	236	LYS
4	R	242	GLU
5	S	4	ASN
5	S	9	THR
5	S	10	VAL
5	S	25	LEU
5	S	29	LYS
5	S	55	LEU
5	S	71	LEU
5	S	92	ASN
5	S	118	ASN
5	S	144	LEU
5	S	188	LEU
5	S	206	THR
5	S	231	LYS
6	T	14	ASP
6	T	68	ARG
6	T	94	SER
6	T	123	ASN
6	T	165	ARG
6	T	172	LEU
6	T	181	GLU
6	T	187	GLU
6	T	203	ASN

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Mol	Chain	Res	Type
6	T	206	LYS
6	T	221	ASN
7	U	26	THR
7	U	28	GLN
7	U	34	LEU
7	U	83	ASN
7	U	115	LEU
7	U	125	MET
7	U	166	GLN
7	U	171	THR
7	U	181	LYS
7	U	235	ARG
7	U	236	LEU
8	V	13	VAL
8	V	34	LEU
8	V	56	THR
8	V	68	LEU
8	V	116	ASP
8	V	127	LEU
8	V	196	ARG
8	V	198	GLU
9	W	37	ASN
9	W	126	ILE
9	W	171	LEU
9	W	182	TRP
9	W	192	ASP
10	X	2	ASP
10	X	23	ARG
10	X	35	THR
10	X	75	LEU
10	X	91	SER
10	X	144	LEU
10	X	163	LEU
10	X	194	ASP
11	Y	4	LEU
11	Y	35	ILE
11	Y	106	ARG
11	Y	118	ASP
11	Y	148	LEU
12	Z	1	GLN
12	Z	23	LEU
12	Z	49	ASN

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Mol	Chain	Res	Type
12	Z	150	LEU
12	Z	167	LYS
13	a	10	SER
13	a	43	ILE
13	a	48	ASN
13	a	70	LEU
13	a	104	ARG
13	a	138	SER
13	a	161	ARG
13	a	187	ARG
13	a	192	SER
13	a	212	LEU
13	a	215	GLU
13	a	223	LYS
14	b	9	LYS
14	b	21	THR
14	b	36	ARG
14	b	119	VAL
14	b	178	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (162) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	30	GLN
1	A	94	HIS
2	B	20	GLN
2	B	58	GLN
2	B	95	GLN
2	B	102	ASN
2	B	119	GLN
2	B	123	GLN
2	B	124	HIS
2	B	155	ASN
2	B	226	GLN
2	B	232	GLN
3	C	17	GLN
3	C	38	ASN
3	C	53	GLN
3	C	77	ASN
3	C	116	GLN
3	C	120	GLN
3	C	147	GLN

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Mol	Chain	Res	Type
3	C	160	GLN
4	D	15	GLN
4	D	91	HIS
4	D	100	ASN
4	D	160	ASN
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	151	ASN
5	E	184	ASN
6	F	19	GLN
6	F	86	ASN
6	F	117	GLN
6	F	123	ASN
7	G	30	ASN
7	G	83	ASN
7	G	114	ASN
7	G	117	GLN
7	G	121	GLN
7	G	166	GLN
7	G	175	ASN
7	G	186	ASN
7	G	212	ASN
7	G	231	ASN
8	H	30	ASN
8	H	57	GLN
8	H	66	HIS
8	H	172	ASN
8	H	189	ASN
9	I	31	GLN
9	I	37	ASN
9	I	63	ASN
10	J	37	GLN
10	J	55	GLN
10	J	63	ASN
10	J	146	HIS
10	J	147	HIS
10	J	191	GLN

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Mol	Chain	Res	Type
11	K	9	GLN
11	K	85	ASN
11	K	143	ASN
11	K	176	ASN
11	K	208	ASN
12	L	3	ASN
12	L	36	ASN
12	L	49	ASN
12	L	70	ASN
12	L	76	HIS
12	L	79	HIS
12	L	80	ASN
12	L	152	ASN
12	L	153	GLN
12	L	158	ASN
12	L	197	GLN
13	M	48	ASN
13	M	108	ASN
13	M	194	ASN
13	M	213	GLN
14	N	38	HIS
14	N	141	ASN
14	N	161	GLN
1	O	30	GLN
1	O	94	HIS
2	P	20	GLN
2	P	58	GLN
2	P	95	GLN
2	P	119	GLN
2	P	123	GLN
2	P	124	HIS
2	P	176	GLN
2	P	226	GLN
2	P	232	GLN
3	Q	17	GLN
3	Q	38	ASN
3	Q	53	GLN
3	Q	77	ASN
3	Q	116	GLN
3	Q	120	GLN
3	Q	147	GLN
3	Q	160	GLN

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Mol	Chain	Res	Type
4	R	15	GLN
4	R	91	HIS
4	R	100	ASN
4	R	149	HIS
4	R	160	ASN
4	R	225	ASN
5	S	68	HIS
5	S	92	ASN
5	S	99	ASN
5	S	116	GLN
5	S	118	ASN
5	S	120	GLN
5	S	151	ASN
5	S	184	ASN
6	T	19	GLN
6	T	86	ASN
6	T	117	GLN
6	T	123	ASN
7	U	30	ASN
7	U	83	ASN
7	U	114	ASN
7	U	117	GLN
7	U	121	GLN
7	U	166	GLN
7	U	175	ASN
7	U	186	ASN
7	U	231	ASN
8	V	30	ASN
8	V	57	GLN
8	V	86	HIS
8	V	172	ASN
8	V	189	ASN
9	W	31	GLN
9	W	63	ASN
9	W	88	GLN
10	X	37	GLN
10	X	55	GLN
10	X	63	ASN
10	X	146	HIS
10	X	191	GLN
11	Y	9	GLN
11	Y	85	ASN

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Mol	Chain	Res	Type
11	Y	143	ASN
11	Y	176	ASN
11	Y	208	ASN
12	Z	3	ASN
12	Z	36	ASN
12	Z	49	ASN
12	Z	70	ASN
12	Z	76	HIS
12	Z	79	HIS
12	Z	80	ASN
12	Z	165	ASN
12	Z	197	GLN
13	a	48	ASN
13	a	108	ASN
14	b	38	HIS
14	b	141	ASN
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 ⓘ

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

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
15	GAU	e	4	14,15	8,8,8	1.23	1 (12%)	8,9,9	1.62	2 (25%)
15	GAU	f	4	8,15	8,8,8	1.21	1 (12%)	8,9,9	1.81	3 (37%)
15	GAU	d	4	15,11	8,8,8	1.38	1 (12%)	8,9,9	1.17	0
15	GAU	c	4	8,15	8,8,8	1.15	1 (12%)	8,9,9	1.33	1 (12%)
15	GAU	h	4	14,15	8,8,8	1.11	1 (12%)	8,9,9	1.04	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
15	GAU	g	4	15,11	8,8,8	1.26	1 (12%)	8,9,9	1.10	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
15	GAU	e	4	14,15	-	0/7/7/7	-
15	GAU	f	4	8,15	-	2/7/7/7	-
15	GAU	d	4	15,11	-	0/7/7/7	-
15	GAU	c	4	8,15	-	0/7/7/7	-
15	GAU	h	4	14,15	-	0/7/7/7	-
15	GAU	g	4	15,11	-	0/7/7/7	-

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
15	e	4	GAU	OE2-CD	-2.68	1.22	1.30
15	h	4	GAU	OE2-CD	-2.54	1.22	1.30
15	d	4	GAU	OE2-CD	-2.43	1.22	1.30
15	g	4	GAU	CG-CD	2.18	1.55	1.50
15	c	4	GAU	OE2-CD	-2.05	1.24	1.30
15	f	4	GAU	CG-CD	2.04	1.55	1.50

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
15	f	4	GAU	CG-CB-CA	-3.50	105.40	112.87
15	e	4	GAU	CG-CB-CA	-2.74	107.03	112.87
15	e	4	GAU	CB-CA-N	-2.67	101.05	109.16
15	c	4	GAU	CB-CA-N	-2.29	102.20	109.16
15	f	4	GAU	OE1-CD-CG	-2.16	116.24	123.09
15	f	4	GAU	CB-CA-C	-2.09	109.39	112.25

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
15	f	4	GAU	C-CA-CB-CG
15	f	4	GAU	N-CA-CB-CG

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
15	e	4	GAU	1	0
15	g	4	GAU	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 7 are monoatomic - leaving 2 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
17	MES	Y	301	-	12,12,12	2.23	1 (8%)	15,16,16	1.13	2 (13%)
17	MES	K	302	-	12,12,12	2.24	1 (8%)	15,16,16	1.16	2 (13%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
17	MES	Y	301	-	-	0/6/14/14	0/1/1/1
17	MES	K	302	-	-	0/6/14/14	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
17	K	302	MES	C8-S	-7.44	1.67	1.77
17	Y	301	MES	C8-S	-7.43	1.67	1.77

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
17	Y	301	MES	O3S-S-C8	2.15	110.22	106.00
17	Y	301	MES	O2S-S-C8	2.13	109.95	106.73
17	K	302	MES	O3S-S-C8	2.09	110.09	106.00
17	K	302	MES	O2S-S-C8	2.06	109.84	106.73

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%)	-0.44	2 (0%) 82 81	43, 59, 89, 130	0
1	O	250/250 (100%)	-0.38	2 (0%) 82 81	46, 67, 102, 141	0
2	B	244/258 (94%)	-0.30	2 (0%) 82 81	45, 65, 107, 150	0
2	P	244/258 (94%)	-0.21	4 (1%) 70 68	49, 69, 114, 160	0
3	C	240/254 (94%)	-0.23	2 (0%) 82 81	46, 70, 129, 157	0
3	Q	240/254 (94%)	-0.04	6 (2%) 58 55	53, 81, 151, 165	0
4	D	235/260 (90%)	-0.30	0 100 100	49, 71, 102, 135	0
4	R	235/260 (90%)	-0.23	0 100 100	51, 72, 105, 136	0
5	E	231/234 (98%)	-0.29	0 100 100	51, 74, 104, 154	0
5	S	231/234 (98%)	-0.13	0 100 100	53, 78, 112, 140	0
6	F	243/288 (84%)	-0.36	1 (0%) 88 87	47, 65, 110, 134	0
6	T	243/288 (84%)	-0.23	4 (1%) 70 68	50, 71, 113, 138	0
7	G	241/252 (95%)	-0.38	0 100 100	43, 61, 95, 138	0
7	U	241/252 (95%)	-0.38	0 100 100	45, 62, 93, 119	0
8	H	222/232 (95%)	-0.52	0 100 100	42, 57, 82, 107	0
8	V	222/232 (95%)	-0.50	0 100 100	29, 59, 85, 116	1 (0%)
9	I	204/205 (99%)	-0.59	0 100 100	42, 56, 82, 106	0
9	W	204/205 (99%)	-0.53	0 100 100	44, 59, 84, 118	0
10	J	195/198 (98%)	-0.48	1 (0%) 87 86	41, 59, 85, 121	0
10	X	195/198 (98%)	-0.52	1 (0%) 87 86	45, 61, 86, 147	0
11	K	212/212 (100%)	-0.48	0 100 100	45, 60, 89, 106	0
11	Y	212/212 (100%)	-0.37	1 (0%) 87 86	47, 63, 98, 123	0
12	L	222/222 (100%)	-0.53	0 100 100	45, 61, 88, 109	0
12	Z	222/222 (100%)	-0.53	1 (0%) 87 86	43, 60, 88, 106	1 (0%)

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
13	M	233/246 (94%)	-0.54	1 (0%) 88 87	41, 59, 77, 91	0
13	a	233/246 (94%)	-0.55	1 (0%) 88 87	43, 57, 76, 90	0
14	N	196/196 (100%)	-0.52	0 100 100	42, 55, 80, 108	0
14	b	196/196 (100%)	-0.52	0 100 100	41, 56, 82, 104	0
15	c	2/5 (40%)	-0.56	0 100 100	48, 48, 48, 53	0
15	d	2/5 (40%)	0.22	0 100 100	60, 60, 60, 61	0
15	e	2/5 (40%)	-0.58	0 100 100	45, 45, 45, 54	0
15	f	2/5 (40%)	-0.40	0 100 100	55, 55, 55, 62	0
15	g	2/5 (40%)	0.01	0 100 100	72, 72, 72, 75	0
15	h	2/5 (40%)	-0.73	0 100 100	53, 53, 53, 57	0
All	All	6348/6644 (95%)	-0.39	29 (0%) 87 86	29, 63, 103, 165	2 (0%)

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
10	J	1	MET	5.3
10	X	1	MET	4.9
1	A	1	MET	4.0
2	P	219	ALA	3.9
1	O	1	MET	3.7
2	B	51	VAL	3.5
3	Q	50	LEU	3.5
3	Q	234	ILE	3.0
13	a	1	THR	3.0
3	C	205	ALA	2.8
6	T	178	HIS	2.8
2	P	51	VAL	2.8
2	P	218	GLY	2.7
3	Q	51	LYS	2.7
6	T	243	ILE	2.5
3	Q	205	ALA	2.5
6	T	205	GLU	2.4
12	Z	108[A]	HIS	2.4
1	O	53	SER	2.3
3	Q	201	VAL	2.3
2	B	219	ALA	2.2
6	T	204	LYS	2.2
3	Q	48	SER	2.2

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Mol	Chain	Res	Type	RSRZ
2	P	93	HIS	2.2
3	C	51	LYS	2.1
1	A	201	GLU	2.1
13	M	1	THR	2.1
11	Y	212	GLY	2.1
6	F	243	ILE	2.0

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

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
15	GAU	h	4	9/9	0.90	0.09	51,52,60,67	0
15	GAU	d	4	9/9	0.94	0.08	41,57,60,61	0
15	GAU	e	4	9/9	0.94	0.08	49,55,61,64	0
15	GAU	f	4	9/9	0.94	0.10	55,58,65,71	0
15	GAU	g	4	9/9	0.94	0.10	64,77,81,81	0
15	GAU	c	4	9/9	0.94	0.09	54,58,65,68	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
17	MES	K	302	12/12	0.88	0.14	91,95,99,100	0
17	MES	Y	301	12/12	0.90	0.12	84,89,98,99	0
16	MG	I	301	1/1	0.92	0.10	69,69,69,69	0
16	MG	H	301	1/1	0.97	0.21	57,57,57,57	0
16	MG	N	201	1/1	0.97	0.04	54,54,54,54	0
16	MG	G	301	1/1	0.99	0.02	57,57,57,57	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
16	MG	K	301	1/1	0.99	0.06	61,61,61,61	0
16	MG	L	301	1/1	0.99	0.07	54,54,54,54	0
16	MG	I	302	1/1	1.00	0.06	55,55,55,55	0

6.5 Other polymers [i](#)

There are no such residues in this entry.