



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 09:09 PM GMT

PDB ID : 4Y8T
Title : Yeast 20S proteasome beta2-H116D mutant in complex with Ac-PAE-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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026982
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20026982

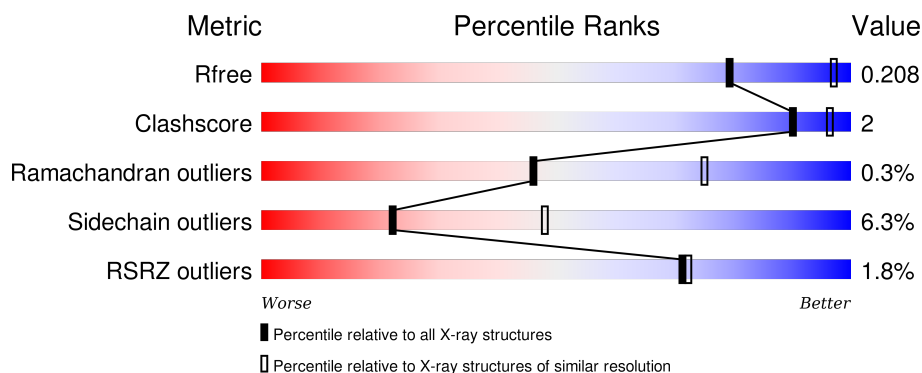
1 Overall quality at a glance

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}	91344	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	250	<div> <div>94%</div> <div>5%</div> </div>
1	O	250	<div> <div>3%</div> <div>94%</div> <div>5%</div> </div>
2	B	258	<div> <div>3%</div> <div>80%</div> <div>13%</div> <div>5%</div> </div>
2	P	258	<div> <div>3%</div> <div>81%</div> <div>13%</div> <div>5%</div> </div>
3	C	254	<div> <div>4%</div> <div>80%</div> <div>13%</div> <div>6%</div> </div>

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Mol	Chain	Length	Quality of chain
3	Q	254	
4	D	260	
4	R	260	
5	E	234	
5	S	234	
6	F	288	
6	T	288	
7	G	252	
7	U	252	
8	H	232	
8	V	232	
9	I	205	
9	W	205	
10	J	198	
10	X	198	
11	K	212	
11	Y	212	
12	L	222	
12	Z	222	
13	M	246	
13	a	246	
14	N	196	
14	b	196	
15	c	5	
15	d	5	

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
16	MG	I	301	-	-	-	X

2 Entry composition

There are 17 unique types of molecules in this entry. The entry contains 50193 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 Proteasome subunit alpha type-7.

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

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

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

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	H	226	Total	C	N	O	S	0	0	0
			1717	1080	296	334	7			
8	V	226	Total	C	N	O	S	0	0	0
			1717	1080	296	334	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	0	0
			1757	1115	303	335	4			

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

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

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

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

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
15	c	5	Total	C	N	O	0	0	0
			28	18	3	7			
15	d	5	Total	C	N	O	0	0	0
			28	18	3	7			

- Molecule 16 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 17 is water.

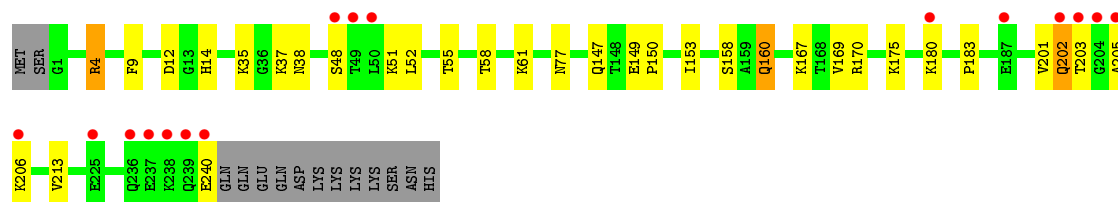
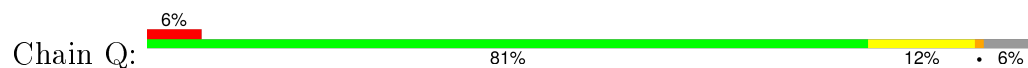
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
17	A	30	Total	O	0	0
			30	30		
17	B	26	Total	O	0	0
			26	26		
17	C	34	Total	O	0	0
			34	34		
17	D	11	Total	O	0	0
			11	11		
17	E	15	Total	O	0	0
			15	15		
17	F	21	Total	O	0	0
			21	21		
17	G	36	Total	O	0	0
			36	36		
17	H	57	Total	O	0	0
			57	57		
17	I	29	Total	O	0	0
			29	29		

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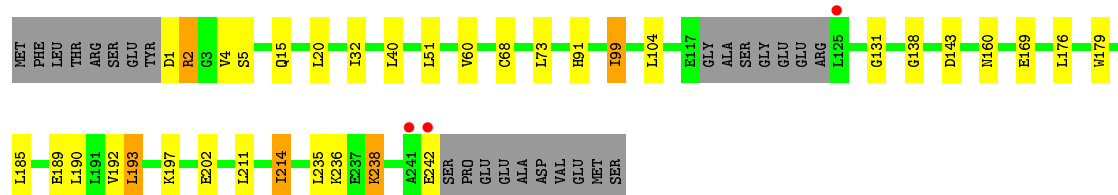
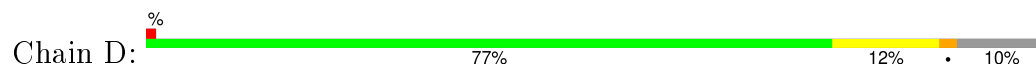
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
17	J	32	Total 32	O 32	0	0
17	K	27	Total 27	O 27	0	0
17	L	39	Total 39	O 39	0	0
17	M	30	Total 30	O 30	0	0
17	N	24	Total 24	O 24	0	0
17	O	18	Total 18	O 18	0	0
17	P	14	Total 14	O 14	0	0
17	Q	15	Total 15	O 15	0	0
17	R	19	Total 19	O 19	0	0
17	S	10	Total 10	O 10	0	0
17	T	24	Total 24	O 24	0	0
17	U	35	Total 35	O 35	0	0
17	V	36	Total 36	O 36	0	0
17	W	23	Total 23	O 23	0	0
17	X	24	Total 24	O 24	0	0
17	Y	36	Total 36	O 36	0	0
17	Z	39	Total 39	O 39	0	0
17	a	36	Total 36	O 36	0	0
17	b	27	Total 27	O 27	0	0
17	c	1	Total 1	O 1	0	0

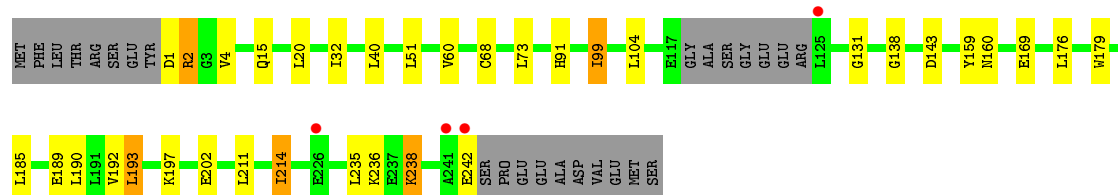
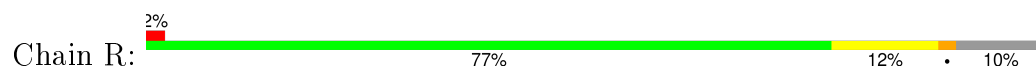
- Molecule 3: Proteasome subunit alpha type-4



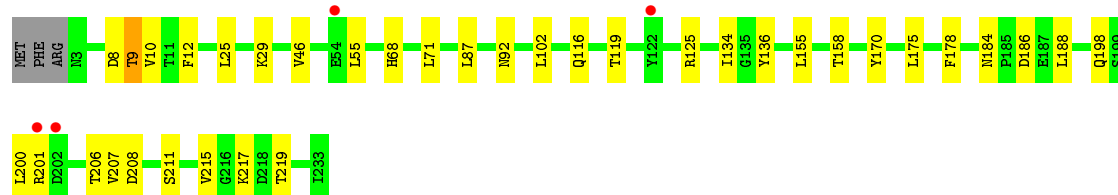
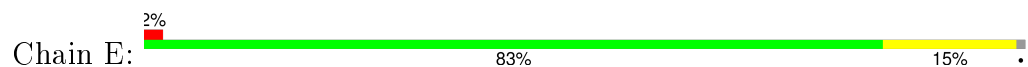
- Molecule 4: Proteasome subunit alpha type-5



- Molecule 4: Proteasome subunit alpha type-5



- Molecule 5: Proteasome subunit alpha type-6



- Molecule 5: Proteasome subunit alpha type-6

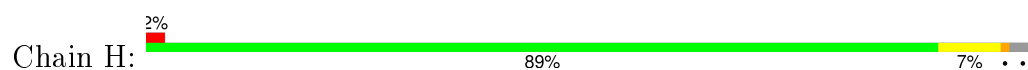
- Molecule 6: Proteasome subunit alpha type-7

- Molecule 6: Proteasome subunit alpha type-7

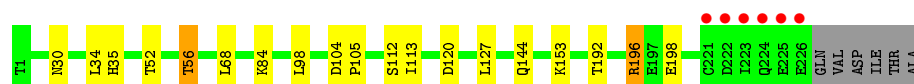
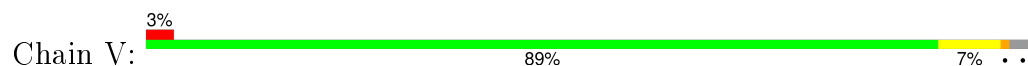
- Molecule 7: Proteasome subunit alpha type-1

- Molecule 7: Proteasome subunit alpha type-1

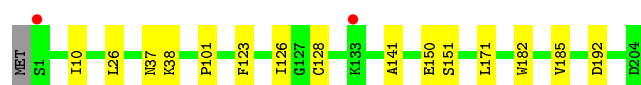
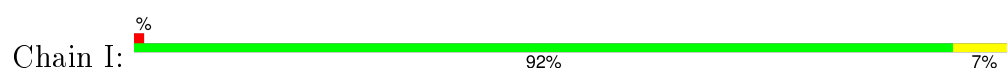
- Molecule 8: Proteasome subunit beta type-2



- Molecule 8: Proteasome subunit beta type-2



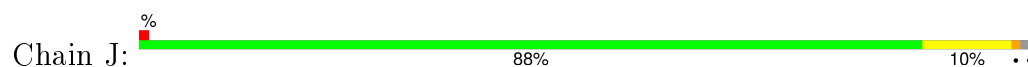
- Molecule 9: Proteasome subunit beta type-3



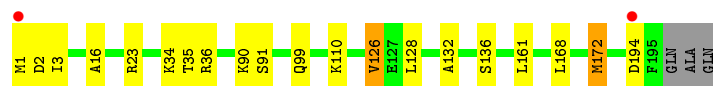
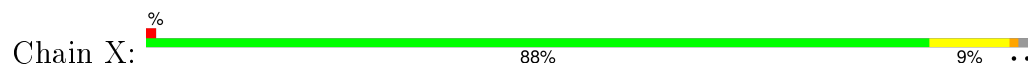
- Molecule 9: Proteasome subunit beta type-3



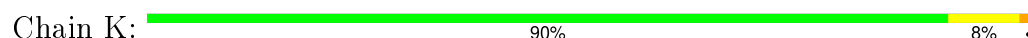
- Molecule 10: Proteasome subunit beta type-4




- Molecule 10: Proteasome subunit beta type-4

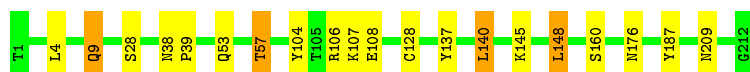


- Molecule 11: Proteasome subunit beta type-5



- Molecule 11: Proteasome subunit beta type-5

Chain Y:  91% 8% •



- Molecule 12: Proteasome subunit beta type-6

Chain L:  93% 7%




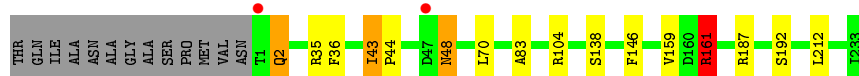
- Molecule 12: Proteasome subunit beta type-6

Chain Z:  92% 7% •




- Molecule 13: Proteasome subunit beta type-7

Chain M:  88% 5% • 5%

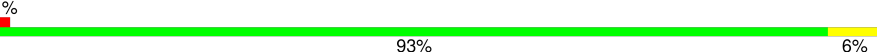


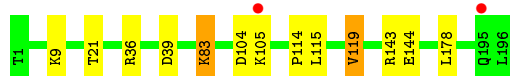
- Molecule 13: Proteasome subunit beta type-7

Chain a:  90% • 5%



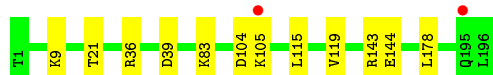
- Molecule 14: Proteasome subunit beta type-1

Chain N:  93% 6% •




- Molecule 14: Proteasome subunit beta type-1

Chain b:  94% 6%




- Molecule 15: Ac-PAE-ep

Chain c:  80% 20%



- Molecule 15: Ac-PAE-ep

Chain d:  80% 20%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	136.40Å 301.59Å 145.63Å 90.00° 113.03° 90.00°	Depositor
Resolution (Å)	15.00 – 2.70 15.00 – 2.70	Depositor EDS
% Data completeness (in resolution range)	97.8 (15.00-2.70) 97.8 (15.00-2.70)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.28 (at 2.69Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.181 , 0.209 0.183 , 0.208	Depositor DCC
R_{free} test set	14351 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	63.4	Xtriage
Anisotropy	0.052	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 46.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	1 of 287021 reflections (0.000%)	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	50193	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.58% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.27	0/1952	0.51	0/2642
1	O	0.27	0/1952	0.51	0/2642
2	B	0.28	0/1934	0.54	0/2618
2	P	0.27	0/1934	0.54	0/2618
3	C	0.28	0/1910	0.55	0/2586
3	Q	0.28	0/1910	0.55	0/2586
4	D	0.28	0/1837	0.60	2/2475 (0.1%)
4	R	0.28	0/1837	0.59	1/2475 (0.0%)
5	E	0.28	0/1800	0.52	0/2433
5	S	0.27	0/1800	0.52	0/2433
6	F	0.28	0/1932	0.50	0/2609
6	T	0.28	0/1932	0.50	0/2609
7	G	0.28	0/1945	0.53	0/2634
7	U	0.28	0/1945	0.52	0/2634
8	H	0.26	0/1747	0.52	0/2369
8	V	0.26	0/1747	0.52	0/2369
9	I	0.28	0/1611	0.51	0/2174
9	W	0.28	0/1611	0.52	0/2174
10	J	0.27	0/1589	0.51	0/2142
10	X	0.27	0/1589	0.51	0/2142
11	K	0.27	0/1681	0.53	0/2274
11	Y	0.27	0/1681	0.53	0/2274
12	L	0.27	0/1795	0.53	0/2420
12	Z	0.27	0/1795	0.52	0/2420
13	M	0.28	0/1855	0.57	1/2514 (0.0%)
13	a	0.27	0/1855	0.57	1/2514 (0.0%)
14	N	0.26	0/1541	0.50	0/2087
14	b	0.27	0/1541	0.49	0/2087
15	c	1.86	1/13 (7.7%)	0.63	0/18
15	d	1.90	1/13 (7.7%)	0.67	0/18
All	All	0.28	2/50284 (0.0%)	0.53	5/67990 (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
2	B	0	1
2	P	0	1
All	All	0	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
15	d	2	PRO	CA-C	-6.15	1.40	1.52
15	c	2	PRO	CA-C	-6.02	1.40	1.52

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	238	LYS	CB-CG-CD	9.55	136.42	111.60
4	R	238	LYS	CB-CG-CD	9.47	136.22	111.60
13	M	161	ARG	NE-CZ-NH1	6.99	123.80	120.30
13	a	161	ARG	NE-CZ-NH1	6.85	123.73	120.30
4	D	238	LYS	CA-CB-CG	6.35	127.37	113.40

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	50	LYS	Peptide
2	P	50	LYS	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	1915	0	1929	6	0
1	O	1915	0	1929	8	0
2	B	1904	0	1904	16	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	P	1904	0	1904	15	0
3	C	1881	0	1895	11	0
3	Q	1881	0	1895	9	0
4	D	1813	0	1797	13	0
4	R	1813	0	1797	13	0
5	E	1773	0	1775	11	0
5	S	1773	0	1775	10	0
6	F	1892	0	1883	8	0
6	T	1892	0	1883	8	0
7	G	1907	0	1901	10	0
7	U	1907	0	1901	9	0
8	H	1717	0	1716	9	0
8	V	1717	0	1716	9	0
9	I	1581	0	1574	6	0
9	W	1581	0	1574	6	0
10	J	1561	0	1569	9	0
10	X	1561	0	1569	8	0
11	K	1644	0	1595	12	0
11	Y	1644	0	1595	10	0
12	L	1757	0	1711	6	0
12	Z	1757	0	1711	7	0
13	M	1824	0	1832	9	0
13	a	1824	0	1832	0	0
14	N	1512	0	1478	4	0
14	b	1512	0	1478	0	0
15	c	28	0	27	0	0
15	d	28	0	27	0	0
16	G	1	0	0	0	0
16	I	1	0	0	0	0
16	K	1	0	0	0	0
16	N	1	0	0	0	0
16	V	1	0	0	0	0
16	Y	1	0	0	0	0
16	Z	1	0	0	0	0
17	A	30	0	0	0	0
17	B	26	0	0	0	0
17	C	34	0	0	0	0
17	D	11	0	0	0	0
17	E	15	0	0	0	0
17	F	21	0	0	0	0
17	G	36	0	0	0	0
17	H	57	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
17	I	29	0	0	0	0
17	J	32	0	0	0	0
17	K	27	0	0	0	0
17	L	39	0	0	0	0
17	M	30	0	0	1	0
17	N	24	0	0	0	0
17	O	18	0	0	0	0
17	P	14	0	0	0	0
17	Q	15	0	0	0	0
17	R	19	0	0	0	0
17	S	10	0	0	0	0
17	T	24	0	0	1	0
17	U	35	0	0	0	0
17	V	36	0	0	0	0
17	W	23	0	0	0	0
17	X	24	0	0	0	0
17	Y	36	0	0	0	0
17	Z	39	0	0	0	0
17	a	36	0	0	0	0
17	b	27	0	0	0	0
17	c	1	0	0	0	0
All	All	50193	0	49172	209	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 (209) 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)
13:M:2:GLN:NE2	17:M:301:HOH:O	2.18	0.75
11:K:53:GLN:O	11:K:57:THR:HG23	1.97	0.64
4:R:99:ILE:HD11	4:R:104:LEU:HB2	1.80	0.63
2:P:217:LYS:O	2:P:219:ALA:N	2.32	0.63
4:D:99:ILE:HD11	4:D:104:LEU:HB2	1.81	0.63
11:Y:53:GLN:O	11:Y:57:THR:HG23	1.98	0.62
2:B:217:LYS:O	2:B:219:ALA:N	2.32	0.62
1:O:38:LYS:NZ	2:P:57:GLU:OE2	2.28	0.62
12:Z:23:LEU:HD13	12:Z:43:VAL:HG13	1.83	0.60
2:B:12:PHE:H	3:C:17:GLN:HE22	1.50	0.60
12:L:23:LEU:HD13	12:L:43:VAL:HG13	1.83	0.59
13:M:161:ARG:HH11	13:M:161:ARG:HG3	1.67	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:160:GLN:HE21	3:C:160:GLN:HA	1.69	0.58
11:Y:145:LYS:HB2	11:Y:148:LEU:HD13	1.87	0.57
4:R:138:GLY:HA2	4:R:214:ILE:HG12	1.87	0.57
3:Q:160:GLN:HA	3:Q:160:GLN:HE21	1.69	0.57
11:K:145:LYS:HB2	11:K:148:LEU:HD13	1.87	0.56
3:C:201:VAL:O	3:C:202:GLN:CB	2.53	0.56
7:G:23:PHE:O	7:G:26:THR:HB	2.05	0.56
4:D:138:GLY:HA2	4:D:214:ILE:HG12	1.87	0.56
7:U:23:PHE:O	7:U:26:THR:HB	2.06	0.56
8:H:52:THR:O	8:H:56:THR:HB	2.07	0.55
3:Q:201:VAL:O	3:Q:202:GLN:CB	2.53	0.55
2:B:146:GLN:HG2	3:C:57:ILE:HG21	1.89	0.55
4:D:32:ILE:HD12	4:D:192:VAL:HG23	1.89	0.55
4:R:32:ILE:HD12	4:R:192:VAL:HG23	1.89	0.54
2:P:47:ALA:HB1	2:P:64:LYS:HD2	1.90	0.54
8:V:84:LYS:HA	8:V:113:ILE:HD11	1.90	0.54
2:B:180:LYS:O	2:B:183:MET:HB2	2.08	0.54
14:N:83:LYS:HG3	14:N:119:VAL:HG22	1.90	0.54
8:V:52:THR:O	8:V:56:THR:HB	2.08	0.53
9:W:10:ILE:HG21	9:W:141:ALA:HB3	1.91	0.53
2:B:47:ALA:HB1	2:B:64:LYS:HD2	1.90	0.53
8:H:84:LYS:HA	8:H:113:ILE:HD11	1.90	0.53
8:H:104:ASP:HB2	8:H:105:PRO:HD2	1.91	0.53
8:V:98:LEU:HB2	8:V:113:ILE:HG22	1.91	0.52
2:P:180:LYS:O	2:P:183:MET:HB2	2.09	0.52
9:I:10:ILE:HG21	9:I:141:ALA:HB3	1.91	0.52
1:A:176:GLU:HG3	2:B:55:LEU:HD22	1.92	0.52
12:L:31:THR:HG23	12:L:36:ASN:HD21	1.73	0.52
12:Z:31:THR:HG23	12:Z:36:ASN:HD21	1.74	0.52
2:B:35:ILE:HD12	2:B:196:LEU:HG	1.91	0.52
8:H:98:LEU:HB2	8:H:113:ILE:HG22	1.91	0.51
8:V:104:ASP:HB2	8:V:105:PRO:HD2	1.91	0.51
2:P:35:ILE:HD12	2:P:196:LEU:HG	1.92	0.51
2:P:3:ARG:HG3	3:Q:4:ARG:CZ	2.41	0.51
3:Q:51:LYS:O	3:Q:52:LEU:HB2	2.11	0.51
2:B:220:ASN:N	2:B:220:ASN:OD1	2.44	0.51
3:C:35:LYS:HG2	3:C:158:SER:O	2.11	0.51
2:P:220:ASN:N	2:P:220:ASN:OD1	2.44	0.51
1:O:23:TYR:CD1	7:U:12:PRO:HA	2.45	0.50
1:A:1:MET:HG3	6:F:122:TYR:CZ	2.46	0.50
3:C:51:LYS:O	3:C:52:LEU:HB2	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:Q:35:LYS:HG2	3:Q:158:SER:O	2.12	0.50
2:B:221:ASP:O	2:B:223:GLU:N	2.45	0.50
13:M:35:ARG:HH12	14:N:114:PRO:HB3	1.76	0.50
13:M:35:ARG:NH1	14:N:114:PRO:HB3	2.27	0.49
10:J:16:ALA:HB2	10:J:161:LEU:HD21	1.93	0.49
10:X:16:ALA:HB2	10:X:161:LEU:HD21	1.93	0.49
11:Y:28:SER:HB2	12:Z:137:ARG:HH22	1.76	0.49
2:P:221:ASP:O	2:P:223:GLU:N	2.45	0.49
12:L:8:ASN:HA	12:L:30:ILE:O	2.13	0.49
13:M:35:ARG:HD3	13:M:36:PHE:CZ	2.48	0.48
11:Y:140:LEU:HD13	11:Y:160:SER:OG	2.14	0.48
9:W:101:PRO:HB3	9:W:126:ILE:HD12	1.96	0.48
12:Z:8:ASN:HA	12:Z:30:ILE:O	2.13	0.48
9:I:101:PRO:HB3	9:I:126:ILE:HD12	1.96	0.48
5:E:175:LEU:HA	5:E:178:PHE:CE2	2.49	0.48
11:K:140:LEU:HD13	11:K:160:SER:OG	2.14	0.48
11:K:128:CYS:HB2	11:K:137:TYR:CE2	2.49	0.48
7:U:187:GLU:HG2	7:U:192:LYS:HB3	1.96	0.48
5:S:175:LEU:HA	5:S:178:PHE:CE2	2.49	0.48
5:E:136:TYR:CE1	5:E:217:LYS:HA	2.49	0.47
7:G:187:GLU:HG2	7:G:192:LYS:HB3	1.96	0.47
11:Y:128:CYS:HB2	11:Y:137:TYR:CE2	2.49	0.47
5:S:155:LEU:HD13	5:S:158:THR:HB	1.97	0.47
2:P:89:THR:HG21	2:P:117:ILE:CD1	2.44	0.47
5:S:136:TYR:CE1	5:S:217:LYS:HA	2.49	0.47
8:H:196:ARG:NH2	9:I:150:GLU:O	2.48	0.47
1:O:1:MET:HG3	6:T:122:TYR:CZ	2.49	0.47
4:R:1:ASP:O	4:R:2:ARG:HB2	2.15	0.47
5:E:155:LEU:HD13	5:E:158:THR:HB	1.97	0.47
11:Y:107:LYS:HG3	11:Y:108:GLU:HG3	1.98	0.46
1:O:222:LEU:HD13	1:O:232:GLY:HA2	1.97	0.46
13:M:161:ARG:CG	13:M:161:ARG:HH11	2.27	0.46
1:O:122:THR:HG22	2:P:128:ARG:HH21	1.80	0.46
13:M:48:ASN:H	13:M:48:ASN:HD22	1.62	0.46
2:B:89:THR:HG21	2:B:117:ILE:CD1	2.45	0.46
4:D:1:ASP:O	4:D:2:ARG:HB2	2.15	0.46
1:A:222:LEU:HD13	1:A:232:GLY:HA2	1.97	0.46
11:K:28:SER:HB2	12:L:137:ARG:HH22	1.80	0.46
6:F:202:ASP:OD1	6:F:202:ASP:N	2.47	0.46
2:P:151:ASN:HB2	2:P:152:PRO:CD	2.47	0.45
4:R:160:ASN:HB3	4:R:179:TRP:CE2	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:160:ASN:HB3	4:D:179:TRP:CE2	2.51	0.45
7:U:63:ILE:HD12	7:U:215:GLU:HG2	1.98	0.45
3:Q:9:PHE:H	4:R:15:GLN:HE22	1.64	0.45
11:K:107:LYS:HG3	11:K:108:GLU:HG3	1.97	0.45
6:T:13:PRO:O	7:U:24:LYS:HD2	2.17	0.45
11:K:176:ASN:ND2	11:K:187:TYR:OH	2.50	0.45
6:F:34:ILE:HG12	6:F:196:ILE:HD11	1.99	0.45
5:E:200:LEU:O	5:E:201:ARG:HG3	2.17	0.45
3:C:9:PHE:H	4:D:15:GLN:HE22	1.65	0.45
5:S:170:TYR:HB2	5:S:198:GLN:HG3	1.99	0.45
14:N:83:LYS:HG3	14:N:119:VAL:CG2	2.47	0.45
6:T:34:ILE:HG12	6:T:196:ILE:HD11	1.99	0.45
7:G:34:LEU:HD23	7:G:34:LEU:C	2.38	0.45
7:G:63:ILE:HD12	7:G:215:GLU:HG2	1.98	0.45
3:C:160:GLN:HE22	3:C:170:ARG:HE	1.65	0.44
3:Q:160:GLN:HE22	3:Q:170:ARG:HE	1.65	0.44
4:R:159:TYR:CE2	5:S:56:SER:HB3	2.51	0.44
8:V:196:ARG:NH2	9:W:150:GLU:O	2.50	0.44
6:T:148:GLU:HG2	17:T:301:HOH:O	2.16	0.44
1:O:12:PHE:H	2:P:20:GLN:HE22	1.63	0.44
5:S:200:LEU:O	5:S:201:ARG:HG3	2.17	0.44
2:B:151:ASN:HB2	2:B:152:PRO:CD	2.47	0.44
11:Y:28:SER:HB2	12:Z:137:ARG:NH2	2.32	0.44
5:S:68:HIS:HE1	5:S:102:LEU:O	2.00	0.44
11:K:209:ASN:O	9:W:38:LYS:NZ	2.51	0.44
10:X:1:MET:HB3	10:X:34:LYS:HE3	1.99	0.44
5:E:68:HIS:HE1	5:E:102:LEU:O	2.00	0.44
5:E:170:TYR:HB2	5:E:198:GLN:HG3	1.99	0.44
6:T:202:ASP:OD1	6:T:202:ASP:N	2.49	0.44
3:C:149:GLU:HB2	3:C:150:PRO:HD2	2.00	0.43
10:X:132:ALA:HB1	10:X:136:SER:HB2	2.00	0.43
5:S:12:PHE:H	6:T:19:GLN:HE22	1.66	0.43
1:A:12:PHE:H	2:B:20:GLN:HE22	1.66	0.43
8:H:35:HIS:HB3	8:H:56:THR:HG21	1.99	0.43
8:H:84:LYS:CA	8:H:113:ILE:HD11	2.48	0.43
10:J:132:ALA:HB1	10:J:136:SER:HB2	2.00	0.43
7:U:34:LEU:C	7:U:34:LEU:HD23	2.39	0.43
7:G:73:VAL:CG1	7:G:133:THR:HB	2.49	0.43
11:Y:176:ASN:ND2	11:Y:187:TYR:OH	2.50	0.43
7:G:195:GLU:HG3	7:G:235:ARG:HG3	2.01	0.43
10:X:126:VAL:HG13	10:X:128:LEU:HG	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:M:159:VAL:HG23	13:M:159:VAL:O	2.19	0.43
6:T:240:GLN:HA	6:T:240:GLN:HE21	1.84	0.43
8:V:35:HIS:HB3	8:V:56:THR:HG21	1.99	0.43
4:R:4:VAL:HG13	4:R:15:GLN:HG3	2.01	0.43
5:E:12:PHE:H	6:F:19:GLN:HE22	1.67	0.43
7:U:195:GLU:HG3	7:U:235:ARG:HG3	2.01	0.43
10:J:1:MET:HB3	10:J:34:LYS:HE3	2.00	0.43
4:D:4:VAL:HG13	4:D:15:GLN:HG3	2.01	0.43
3:Q:149:GLU:HB2	3:Q:150:PRO:HD2	2.00	0.43
1:A:122:THR:CG2	2:B:128:ARG:HH21	2.32	0.42
9:I:26:LEU:HD21	9:I:185:VAL:HG23	2.00	0.42
4:R:73:LEU:HD12	4:R:131:GLY:HA3	2.01	0.42
4:D:91:HIS:HB3	4:D:99:ILE:HG21	2.02	0.42
1:O:1:MET:CG	1:O:2:THR:N	2.82	0.42
12:L:195:HIS:HD2	12:L:197:GLN:H	1.66	0.42
4:D:185:LEU:O	4:D:189:GLU:HG3	2.19	0.42
1:O:122:THR:CG2	2:P:128:ARG:HH21	2.32	0.42
9:W:123:PHE:HA	9:W:128:CYS:O	2.19	0.42
13:M:43:ILE:HA	13:M:44:PRO:HD3	1.86	0.42
12:Z:195:HIS:HD2	12:Z:197:GLN:H	1.67	0.42
9:W:26:LEU:HD21	9:W:185:VAL:HG23	2.00	0.42
4:R:185:LEU:O	4:R:189:GLU:HG3	2.20	0.42
4:D:73:LEU:HD12	4:D:131:GLY:HA3	2.02	0.42
6:F:13:PRO:O	7:G:24:LYS:HD2	2.20	0.42
10:J:126:VAL:HG13	10:J:128:LEU:HG	2.01	0.42
8:V:84:LYS:CA	8:V:113:ILE:HD11	2.49	0.42
7:G:73:VAL:HG12	7:G:133:THR:HB	2.02	0.42
3:C:12:ASP:OD2	3:C:14:HIS:ND1	2.52	0.42
7:U:73:VAL:CG1	7:U:133:THR:HB	2.50	0.42
6:T:146:MET:CE	6:T:161:THR:HB	2.50	0.42
9:I:123:PHE:HA	9:I:128:CYS:O	2.19	0.42
4:D:91:HIS:HB3	4:D:99:ILE:CG2	2.50	0.42
4:D:5:SER:HB3	5:E:125:ARG:HD3	2.02	0.42
10:J:126:VAL:CG1	10:J:128:LEU:HG	2.49	0.42
5:E:134:ILE:HD12	5:E:215:VAL:HG12	2.02	0.42
6:F:240:GLN:HA	6:F:240:GLN:HE21	1.84	0.41
2:B:50:LYS:HA	2:B:50:LYS:HD3	1.52	0.41
10:X:126:VAL:CG1	10:X:128:LEU:HG	2.50	0.41
5:S:134:ILE:HD12	5:S:215:VAL:HG12	2.02	0.41
10:J:36:ARG:HD3	10:J:36:ARG:HA	1.92	0.41
8:V:192:THR:HG22	8:V:192:THR:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:159:TRP:HH2	3:C:50:LEU:HD22	1.85	0.41
8:V:112:SER:OG	8:V:120:ASP:OD1	2.27	0.41
3:Q:12:ASP:OD2	3:Q:14:HIS:ND1	2.53	0.41
2:P:139:TYR:CD1	2:P:224:VAL:HG21	2.56	0.41
6:F:146:MET:CE	6:F:161:THR:HB	2.50	0.41
4:R:91:HIS:HB3	4:R:99:ILE:HG21	2.01	0.41
5:E:9:THR:HG21	5:E:119:THR:HA	2.01	0.41
10:X:36:ARG:HD3	10:X:36:ARG:HA	1.93	0.41
7:G:78:ILE:N	7:G:79:PRO:CD	2.84	0.41
2:B:139:TYR:CD1	2:B:224:VAL:HG21	2.55	0.41
1:A:1:MET:CG	1:A:2:THR:N	2.82	0.41
11:Y:38:ASN:HB2	11:Y:39:PRO:CD	2.51	0.41
10:X:168:LEU:O	10:X:172:MET:HB2	2.21	0.41
7:G:227:LEU:HB3	7:G:231:ASN:HB2	2.02	0.41
11:K:28:SER:HB2	12:L:137:ARG:NH2	2.36	0.41
9:I:38:LYS:NZ	11:Y:209:ASN:O	2.53	0.41
2:P:50:LYS:HD3	2:P:50:LYS:HA	1.48	0.41
8:H:192:THR:HG22	8:H:192:THR:O	2.21	0.41
11:K:38:ASN:HB2	11:K:39:PRO:CD	2.51	0.41
8:H:112:SER:OG	8:H:120:ASP:OD1	2.26	0.41
11:K:4:LEU:HD13	11:K:161:ILE:HD11	2.03	0.41
4:R:91:HIS:HB3	4:R:99:ILE:CG2	2.50	0.41
10:J:50:ALA:O	11:K:91:LYS:NZ	2.54	0.41
5:S:9:THR:HG21	5:S:119:THR:HA	2.01	0.41
4:D:193:LEU:HD22	4:D:211:LEU:HD11	2.03	0.41
10:J:194:ASP:N	10:J:194:ASP:OD1	2.55	0.40
10:J:168:LEU:O	10:J:172:MET:HB2	2.22	0.40
10:X:194:ASP:N	10:X:194:ASP:OD1	2.54	0.40
6:F:238:PHE:O	6:F:242:GLU:HG2	2.22	0.40
5:E:87:LEU:HD23	5:E:87:LEU:HA	1.95	0.40
7:U:78:ILE:N	7:U:79:PRO:CD	2.84	0.40
12:Z:146:ILE:HG22	12:Z:150:LEU:HD22	2.03	0.40
4:R:193:LEU:HD22	4:R:211:LEU:HD11	2.03	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%)	241 (97%)	5 (2%)	2 (1%)	24	51
1	O	248/250 (99%)	240 (97%)	6 (2%)	2 (1%)	24	51
2	B	242/258 (94%)	233 (96%)	7 (3%)	2 (1%)	24	51
2	P	242/258 (94%)	233 (96%)	7 (3%)	2 (1%)	24	51
3	C	238/254 (94%)	232 (98%)	3 (1%)	3 (1%)	15	37
3	Q	238/254 (94%)	232 (98%)	3 (1%)	3 (1%)	15	37
4	D	231/260 (89%)	228 (99%)	2 (1%)	1 (0%)	39	69
4	R	231/260 (89%)	228 (99%)	2 (1%)	1 (0%)	39	69
5	E	229/234 (98%)	219 (96%)	10 (4%)	0	100	100
5	S	229/234 (98%)	220 (96%)	9 (4%)	0	100	100
6	F	241/288 (84%)	232 (96%)	9 (4%)	0	100	100
6	T	241/288 (84%)	232 (96%)	9 (4%)	0	100	100
7	G	239/252 (95%)	236 (99%)	3 (1%)	0	100	100
7	U	239/252 (95%)	236 (99%)	3 (1%)	0	100	100
8	H	224/232 (97%)	218 (97%)	6 (3%)	0	100	100
8	V	224/232 (97%)	218 (97%)	6 (3%)	0	100	100
9	I	202/205 (98%)	197 (98%)	5 (2%)	0	100	100
9	W	202/205 (98%)	197 (98%)	5 (2%)	0	100	100
10	J	193/198 (98%)	186 (96%)	7 (4%)	0	100	100
10	X	193/198 (98%)	186 (96%)	7 (4%)	0	100	100
11	K	210/212 (99%)	204 (97%)	5 (2%)	1 (0%)	34	63
11	Y	210/212 (99%)	204 (97%)	5 (2%)	1 (0%)	34	63
12	L	220/222 (99%)	215 (98%)	5 (2%)	0	100	100
12	Z	220/222 (99%)	215 (98%)	5 (2%)	0	100	100
13	M	231/246 (94%)	223 (96%)	7 (3%)	1 (0%)	39	69

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
13	a	231/246 (94%)	223 (96%)	7 (3%)	1 (0%)	39	69
14	N	194/196 (99%)	190 (98%)	4 (2%)	0	100	100
14	b	194/196 (99%)	190 (98%)	4 (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
All	All	6288/6624 (95%)	6112 (97%)	156 (2%)	20 (0%)	46	75

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	2	THR
2	B	218	GLY
2	B	222	GLY
3	C	202	GLN
3	C	205	ALA
1	O	2	THR
2	P	218	GLY
2	P	222	GLY
3	Q	202	GLN
3	Q	205	ALA
4	D	2	ARG
4	R	2	ARG
3	C	183	PRO
3	Q	183	PRO
11	K	9	GLN
11	Y	9	GLN
1	A	166	LYS
1	O	166	LYS
13	M	83	ALA
13	a	83	ALA

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%)	201 (96%)	8 (4%)	40	71
1	O	209/209 (100%)	201 (96%)	8 (4%)	40	71
2	B	203/216 (94%)	188 (93%)	15 (7%)	17	39
2	P	203/216 (94%)	188 (93%)	15 (7%)	17	39
3	C	212/226 (94%)	193 (91%)	19 (9%)	12	27
3	Q	212/226 (94%)	193 (91%)	19 (9%)	12	27
4	D	194/215 (90%)	176 (91%)	18 (9%)	11	25
4	R	194/215 (90%)	176 (91%)	18 (9%)	11	25
5	E	190/193 (98%)	172 (90%)	18 (10%)	11	24
5	S	190/193 (98%)	173 (91%)	17 (9%)	12	27
6	F	201/239 (84%)	186 (92%)	15 (8%)	17	38
6	T	201/239 (84%)	186 (92%)	15 (8%)	17	38
7	G	206/210 (98%)	192 (93%)	14 (7%)	20	43
7	U	206/210 (98%)	194 (94%)	12 (6%)	25	52
8	H	185/190 (97%)	176 (95%)	9 (5%)	31	61
8	V	185/190 (97%)	176 (95%)	9 (5%)	31	61
9	I	172/173 (99%)	167 (97%)	5 (3%)	50	80
9	W	172/173 (99%)	167 (97%)	5 (3%)	50	80
10	J	173/175 (99%)	163 (94%)	10 (6%)	25	52
10	X	173/175 (99%)	163 (94%)	10 (6%)	25	52
11	K	169/169 (100%)	162 (96%)	7 (4%)	37	69
11	Y	169/169 (100%)	162 (96%)	7 (4%)	37	69
12	L	185/185 (100%)	177 (96%)	8 (4%)	35	66
12	Z	185/185 (100%)	176 (95%)	9 (5%)	31	61
13	M	199/208 (96%)	188 (94%)	11 (6%)	27	55
13	a	199/208 (96%)	188 (94%)	11 (6%)	27	55
14	N	162/162 (100%)	150 (93%)	12 (7%)	17	39
14	b	162/162 (100%)	150 (93%)	12 (7%)	17	39
15	c	1/1 (100%)	1 (100%)	0	100	100
15	d	1/1 (100%)	1 (100%)	0	100	100
All	All	5322/5542 (96%)	4986 (94%)	336 (6%)	22	48

All (336) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	THR
1	A	17	LYS
1	A	58	SER
1	A	59	GLU
1	A	61	LEU
1	A	122	THR
1	A	157	PHE
1	A	231	LYS
2	B	50	LYS
2	B	52	THR
2	B	54	THR
2	B	55	LEU
2	B	58	GLN
2	B	65	LEU
2	B	67	LYS
2	B	102	ASN
2	B	119	GLN
2	B	180	LYS
2	B	191	LEU
2	B	194	LYS
2	B	220	ASN
2	B	238	LEU
2	B	244	THR
3	C	4	ARG
3	C	37	LYS
3	C	38	ASN
3	C	48	SER
3	C	55	THR
3	C	58	THR
3	C	61	LYS
3	C	77	ASN
3	C	147	GLN
3	C	153	ILE
3	C	160	GLN
3	C	167	LYS
3	C	169	VAL
3	C	175	LYS
3	C	180	LYS
3	C	203	THR
3	C	206	LYS
3	C	213	VAL
3	C	240	GLU
4	D	20	LEU

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Mol	Chain	Res	Type
4	D	40	LEU
4	D	51	LEU
4	D	60	VAL
4	D	68	CYS
4	D	99	ILE
4	D	143	ASP
4	D	169	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	238	LYS
4	D	242	GLU
5	E	8	ASP
5	E	9	THR
5	E	10	VAL
5	E	25	LEU
5	E	29	LYS
5	E	46	VAL
5	E	55	LEU
5	E	71	LEU
5	E	92	ASN
5	E	116	GLN
5	E	184	ASN
5	E	186	ASP
5	E	188	LEU
5	E	206	THR
5	E	207	VAL
5	E	208	ASP
5	E	211	SER
5	E	219	THR
6	F	14	ASP
6	F	58	GLN
6	F	94	SER
6	F	117	GLN
6	F	123	ASN
6	F	139	LYS
6	F	165	ARG

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Mol	Chain	Res	Type
6	F	172	LEU
6	F	181	GLU
6	F	201	GLU
6	F	202	ASP
6	F	203	ASN
6	F	214	TRP
6	F	221	ASN
6	F	228	LYS
7	G	13	GLU
7	G	26	THR
7	G	28	GLN
7	G	34	LEU
7	G	68	ARG
7	G	83	ASN
7	G	115	LEU
7	G	117	GLN
7	G	122	ARG
7	G	125	MET
7	G	207	THR
7	G	208	GLU
7	G	235	ARG
7	G	236	LEU
8	H	30	ASN
8	H	34	LEU
8	H	56	THR
8	H	68	LEU
8	H	127	LEU
8	H	144	GLN
8	H	153	LYS
8	H	196	ARG
8	H	198	GLU
9	I	37	ASN
9	I	151	SER
9	I	171	LEU
9	I	182	TRP
9	I	192	ASP
10	J	2	ASP
10	J	3	ILE
10	J	23	ARG
10	J	35	THR
10	J	90	LYS
10	J	91	SER

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Mol	Chain	Res	Type
10	J	99	GLN
10	J	110	LYS
10	J	126	VAL
10	J	172	MET
11	K	4	LEU
11	K	9	GLN
11	K	57	THR
11	K	104	TYR
11	K	106	ARG
11	K	140	LEU
11	K	148	LEU
12	L	3	ASN
12	L	23	LEU
12	L	49	ASN
12	L	124	SER
12	L	132	GLU
12	L	150	LEU
12	L	161	GLU
12	L	173	LYS
13	M	2	GLN
13	M	43	ILE
13	M	48	ASN
13	M	70	LEU
13	M	104	ARG
13	M	138	SER
13	M	146	PHE
13	M	161	ARG
13	M	187	ARG
13	M	192	SER
13	M	212	LEU
14	N	9	LYS
14	N	21	THR
14	N	36	ARG
14	N	39	ASP
14	N	83	LYS
14	N	104	ASP
14	N	105	LYS
14	N	115	LEU
14	N	119	VAL
14	N	143	ARG
14	N	144	GLU
14	N	178	LEU

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Mol	Chain	Res	Type
1	O	2	THR
1	O	17	LYS
1	O	58	SER
1	O	59	GLU
1	O	61	LEU
1	O	122	THR
1	O	157	PHE
1	O	231	LYS
2	P	50	LYS
2	P	52	THR
2	P	54	THR
2	P	55	LEU
2	P	58	GLN
2	P	65	LEU
2	P	67	LYS
2	P	102	ASN
2	P	119	GLN
2	P	180	LYS
2	P	191	LEU
2	P	194	LYS
2	P	220	ASN
2	P	238	LEU
2	P	244	THR
3	Q	4	ARG
3	Q	37	LYS
3	Q	38	ASN
3	Q	48	SER
3	Q	55	THR
3	Q	58	THR
3	Q	61	LYS
3	Q	77	ASN
3	Q	147	GLN
3	Q	153	ILE
3	Q	160	GLN
3	Q	167	LYS
3	Q	169	VAL
3	Q	175	LYS
3	Q	180	LYS
3	Q	203	THR
3	Q	206	LYS
3	Q	213	VAL
3	Q	240	GLU

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Mol	Chain	Res	Type
4	R	20	LEU
4	R	40	LEU
4	R	51	LEU
4	R	60	VAL
4	R	68	CYS
4	R	99	ILE
4	R	143	ASP
4	R	169	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	238	LYS
4	R	242	GLU
5	S	8	ASP
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	116	GLN
5	S	184	ASN
5	S	186	ASP
5	S	188	LEU
5	S	206	THR
5	S	207	VAL
5	S	208	ASP
5	S	211	SER
5	S	219	THR
6	T	14	ASP
6	T	58	GLN
6	T	94	SER
6	T	117	GLN
6	T	123	ASN
6	T	139	LYS
6	T	165	ARG

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Mol	Chain	Res	Type
6	T	172	LEU
6	T	181	GLU
6	T	201	GLU
6	T	202	ASP
6	T	203	ASN
6	T	214	TRP
6	T	221	ASN
6	T	228	LYS
7	U	13	GLU
7	U	26	THR
7	U	28	GLN
7	U	68	ARG
7	U	83	ASN
7	U	115	LEU
7	U	117	GLN
7	U	122	ARG
7	U	125	MET
7	U	208	GLU
7	U	235	ARG
7	U	236	LEU
8	V	30	ASN
8	V	34	LEU
8	V	56	THR
8	V	68	LEU
8	V	127	LEU
8	V	144	GLN
8	V	153	LYS
8	V	196	ARG
8	V	198	GLU
9	W	37	ASN
9	W	151	SER
9	W	171	LEU
9	W	182	TRP
9	W	192	ASP
10	X	2	ASP
10	X	3	ILE
10	X	23	ARG
10	X	35	THR
10	X	90	LYS
10	X	91	SER
10	X	99	GLN
10	X	110	LYS

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Mol	Chain	Res	Type
10	X	126	VAL
10	X	172	MET
11	Y	4	LEU
11	Y	9	GLN
11	Y	57	THR
11	Y	104	TYR
11	Y	106	ARG
11	Y	140	LEU
11	Y	148	LEU
12	Z	3	ASN
12	Z	23	LEU
12	Z	49	ASN
12	Z	124	SER
12	Z	132	GLU
12	Z	150	LEU
12	Z	161	GLU
12	Z	172	LEU
12	Z	173	LYS
13	a	2	GLN
13	a	43	ILE
13	a	48	ASN
13	a	70	LEU
13	a	104	ARG
13	a	138	SER
13	a	146	PHE
13	a	161	ARG
13	a	187	ARG
13	a	192	SER
13	a	212	LEU
14	b	9	LYS
14	b	21	THR
14	b	36	ARG
14	b	39	ASP
14	b	83	LYS
14	b	104	ASP
14	b	105	LYS
14	b	115	LEU
14	b	119	VAL
14	b	143	ARG
14	b	144	GLU
14	b	178	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (126) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	94	HIS
2	B	20	GLN
2	B	58	GLN
2	B	95	GLN
2	B	119	GLN
2	B	123	GLN
2	B	155	ASN
3	C	17	GLN
3	C	38	ASN
3	C	77	ASN
3	C	116	GLN
3	C	120	GLN
3	C	147	GLN
3	C	160	GLN
3	C	233	GLN
4	D	15	GLN
4	D	91	HIS
4	D	100	ASN
4	D	146	GLN
4	D	225	ASN
5	E	68	HIS
5	E	92	ASN
5	E	99	ASN
5	E	116	GLN
5	E	118	ASN
5	E	120	GLN
5	E	147	GLN
5	E	184	ASN
6	F	19	GLN
6	F	86	ASN
6	F	117	GLN
6	F	123	ASN
6	F	191	GLN
6	F	240	GLN
7	G	30	ASN
7	G	83	ASN
7	G	114	ASN
7	G	117	GLN
7	G	121	GLN
7	G	167	GLN
7	G	175	ASN
8	H	66	HIS

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Mol	Chain	Res	Type
8	H	114	HIS
8	H	165	ASN
9	I	37	ASN
10	J	55	GLN
10	J	147	HIS
11	K	85	ASN
11	K	176	ASN
11	K	190	ASN
11	K	208	ASN
12	L	3	ASN
12	L	49	ASN
12	L	70	ASN
12	L	80	ASN
12	L	158	ASN
13	M	18	ASN
13	M	48	ASN
13	M	102	GLN
13	M	179	ASN
13	M	194	ASN
13	M	213	GLN
14	N	161	GLN
1	O	94	HIS
2	P	20	GLN
2	P	58	GLN
2	P	95	GLN
2	P	119	GLN
2	P	123	GLN
3	Q	17	GLN
3	Q	38	ASN
3	Q	77	ASN
3	Q	116	GLN
3	Q	120	GLN
3	Q	147	GLN
3	Q	160	GLN
3	Q	233	GLN
4	R	15	GLN
4	R	91	HIS
4	R	100	ASN
4	R	146	GLN
4	R	225	ASN
5	S	68	HIS
5	S	92	ASN

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Mol	Chain	Res	Type
5	S	99	ASN
5	S	116	GLN
5	S	118	ASN
5	S	120	GLN
5	S	184	ASN
6	T	19	GLN
6	T	86	ASN
6	T	117	GLN
6	T	123	ASN
6	T	191	GLN
6	T	240	GLN
7	U	30	ASN
7	U	83	ASN
7	U	114	ASN
7	U	117	GLN
7	U	121	GLN
7	U	167	GLN
7	U	175	ASN
7	U	186	ASN
8	V	66	HIS
8	V	114	HIS
8	V	165	ASN
9	W	37	ASN
10	X	55	GLN
10	X	86	GLN
10	X	147	HIS
11	Y	85	ASN
11	Y	176	ASN
11	Y	190	ASN
11	Y	208	ASN
12	Z	3	ASN
12	Z	49	ASN
12	Z	70	ASN
12	Z	80	ASN
12	Z	158	ASN
13	a	18	ASN
13	a	48	ASN
13	a	102	GLN
13	a	179	ASN
13	a	194	ASN
13	a	213	GLN
14	b	161	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 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	c	4	15,14	5,8,8	1.41	1 (20%)	4,9,9	2.02	1 (25%)
15	GAU	d	4	15,14	5,8,8	1.39	1 (20%)	4,9,9	2.01	1 (25%)

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	c	4	15,14	-	0/5/7/7	0/0/0/0
15	GAU	d	4	15,14	-	0/5/7/7	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
15	d	4	GAU	C-CA	3.07	1.56	1.52
15	c	4	GAU	C-CA	3.12	1.56	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
15	c	4	GAU	CG-CB-CA	-3.84	104.22	113.80
15	d	4	GAU	CG-CB-CA	-3.83	104.26	113.80

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 7 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

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%)	-0.47	3 (1%) 81 81	40, 61, 94, 139	0
1	O	250/250 (100%)	-0.45	7 (2%) 56 57	47, 69, 110, 138	0
2	B	244/258 (94%)	-0.42	7 (2%) 55 55	44, 64, 112, 159	0
2	P	244/258 (94%)	-0.31	8 (3%) 50 50	51, 71, 114, 163	0
3	C	240/254 (94%)	-0.21	11 (4%) 36 35	49, 71, 131, 157	0
3	Q	240/254 (94%)	-0.06	16 (6%) 21 19	54, 81, 148, 164	0
4	D	235/260 (90%)	-0.48	3 (1%) 79 79	50, 71, 105, 134	0
4	R	235/260 (90%)	-0.40	4 (1%) 73 74	50, 72, 107, 140	0
5	E	231/234 (98%)	-0.35	4 (1%) 73 74	53, 75, 109, 153	0
5	S	231/234 (98%)	-0.26	4 (1%) 73 74	54, 81, 115, 158	0
6	F	243/288 (84%)	-0.44	7 (2%) 55 55	47, 69, 113, 141	0
6	T	243/288 (84%)	-0.40	5 (2%) 67 68	52, 75, 123, 151	0
7	G	241/252 (95%)	-0.49	5 (2%) 67 68	47, 62, 95, 146	0
7	U	241/252 (95%)	-0.41	3 (1%) 81 81	50, 67, 100, 129	0
8	H	226/232 (97%)	-0.52	5 (2%) 65 66	45, 58, 90, 144	0
8	V	226/232 (97%)	-0.45	6 (2%) 58 58	47, 63, 95, 163	0
9	I	204/205 (99%)	-0.71	2 (0%) 84 85	43, 58, 84, 111	0
9	W	204/205 (99%)	-0.67	1 (0%) 91 93	47, 61, 92, 124	0
10	J	195/198 (98%)	-0.55	2 (1%) 84 85	47, 61, 89, 144	0
10	X	195/198 (98%)	-0.49	2 (1%) 84 85	47, 63, 88, 150	0
11	K	212/212 (100%)	-0.58	0 100 100	47, 63, 86, 104	0
11	Y	212/212 (100%)	-0.59	0 100 100	47, 62, 85, 107	0
12	L	222/222 (100%)	-0.58	1 (0%) 91 93	47, 62, 94, 139	0
12	Z	222/222 (100%)	-0.60	1 (0%) 91 93	46, 60, 92, 134	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
13	M	233/246 (94%)	-0.63	2 (0%) 85 86	45, 61, 82, 108	0
13	a	233/246 (94%)	-0.61	1 (0%) 93 94	46, 60, 81, 96	0
14	N	196/196 (100%)	-0.69	2 (1%) 84 85	45, 57, 83, 115	0
14	b	196/196 (100%)	-0.64	2 (1%) 84 85	42, 58, 84, 119	0
15	c	2/5 (40%)	-0.85	0 100 100	53, 53, 53, 55	0
15	d	2/5 (40%)	-1.02	0 100 100	52, 52, 52, 54	0
All	All	6348/6624 (95%)	-0.48	114 (1%) 71 72	40, 65, 107, 164	0

All (114) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	Q	49	THR	7.2
3	Q	236	GLN	6.1
3	Q	50	LEU	6.0
8	V	224	GLN	5.8
3	C	206	LYS	5.5
10	X	1	MET	5.3
3	C	49	THR	5.0
3	C	238	LYS	5.0
9	W	1	SER	4.9
8	V	226	GLU	4.7
5	S	202	ASP	4.7
6	T	243	ILE	4.4
8	V	221	CYS	4.3
3	Q	206	LYS	4.2
5	E	202	ASP	4.2
8	V	222	ASP	4.2
2	B	221	ASP	4.1
2	P	51	VAL	4.1
3	C	236	GLN	4.0
1	A	1	MET	4.0
7	G	2	GLY	3.7
3	Q	202	GLN	3.6
8	H	224	GLN	3.5
2	B	51	VAL	3.5
2	P	59	ASP	3.5
2	P	221	ASP	3.5
10	J	1	MET	3.4
10	X	194	ASP	3.3
4	R	125	LEU	3.3

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Mol	Chain	Res	Type	RSRZ
12	Z	174	TYR	3.2
2	P	222	GLY	3.2
3	Q	239	GLN	3.2
12	L	174	TYR	3.2
6	F	205	GLU	3.2
6	F	202	ASP	3.1
8	H	226	GLU	3.0
1	O	249	ALA	3.0
2	P	218	GLY	3.0
1	A	2	THR	3.0
1	O	250	LEU	3.0
2	B	218	GLY	3.0
13	a	1	THR	2.9
3	C	239	GLN	2.9
6	T	244	ASN	2.8
3	C	205	ALA	2.8
8	H	221	CYS	2.8
9	I	1	SER	2.8
3	C	202	GLN	2.8
10	J	194	ASP	2.8
4	D	125	LEU	2.7
6	T	166	GLN	2.7
2	P	50	LYS	2.7
3	C	175	LYS	2.7
7	G	3	TYR	2.7
2	P	60	THR	2.7
8	V	225	GLU	2.7
3	Q	203	THR	2.7
7	U	2	GLY	2.7
5	E	122	TYR	2.7
5	E	201	ARG	2.7
5	E	54	GLU	2.7
6	F	241	LYS	2.7
4	R	241	ALA	2.6
8	H	222	ASP	2.6
2	B	219	ALA	2.6
14	N	105	LYS	2.6
4	D	242	GLU	2.6
3	Q	238	LYS	2.6
7	U	242	GLN	2.6
3	Q	240	GLU	2.6
4	R	242	GLU	2.5

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Mol	Chain	Res	Type	RSRZ
2	P	219	ALA	2.5
8	V	223	ILE	2.5
3	Q	48	SER	2.5
7	U	188	GLU	2.5
6	F	178	HIS	2.5
1	A	249	ALA	2.5
2	B	220	ASN	2.5
7	G	240	ALA	2.5
7	G	179	LYS	2.5
5	S	201	ARG	2.5
1	O	50	LYS	2.4
6	F	244	ASN	2.4
3	Q	180	LYS	2.4
6	T	241	LYS	2.4
3	C	216	ASP	2.3
3	C	180	LYS	2.3
5	S	180	LYS	2.3
9	I	133	LYS	2.3
3	Q	204	GLY	2.3
3	Q	225	GLU	2.3
13	M	1	THR	2.2
14	b	195	GLN	2.2
14	b	105	LYS	2.2
2	B	242	GLY	2.2
3	C	240	GLU	2.2
6	F	243	ILE	2.2
1	O	231	LYS	2.2
3	Q	187	GLU	2.2
13	M	47	ASP	2.2
8	H	223	ILE	2.2
3	Q	237	GLU	2.2
6	F	181	GLU	2.1
3	Q	205	ALA	2.1
2	B	59	ASP	2.1
4	D	241	ALA	2.1
1	O	2	THR	2.1
6	T	180	PRO	2.1
7	G	242	GLN	2.1
4	R	226	GLU	2.1
5	S	54	GLU	2.1
14	N	195	GLN	2.0
1	O	1	MET	2.0

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Mol	Chain	Res	Type	RSRZ
1	O	201	GLU	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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
15	GAU	c	4	9/9	0.98	0.08	-	58,62,64,64	0
15	GAU	d	4	9/9	0.97	0.11	-	57,59,62,67	0

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
16	MG	I	301	1/1	0.96	0.24	4.88	63,63,63,63	0
16	MG	Z	301	1/1	0.96	0.16	1.31	60,60,60,60	0
16	MG	G	301	1/1	0.96	0.11	-0.18	65,65,65,65	0
16	MG	K	301	1/1	0.97	0.11	-0.27	65,65,65,65	0
16	MG	V	301	1/1	0.97	0.10	-0.71	74,74,74,74	0
16	MG	Y	301	1/1	0.98	0.07	-2.47	56,56,56,56	0
16	MG	N	201	1/1	0.98	0.06	-2.67	57,57,57,57	0

6.5 Other polymers [i](#)

There are no such residues in this entry.