



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:49 PM UTC

PDB ID : 6L9Z / pdb_00006l9z
Title : 338 bp di-nucleosome assembled with linker histone H1.X
Authors : Adhireksan, Z.; Sharma, D.; Lee, P.L.; Davey, C.A.
Deposited on : 2019-11-11
Resolution : 2.50 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<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
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

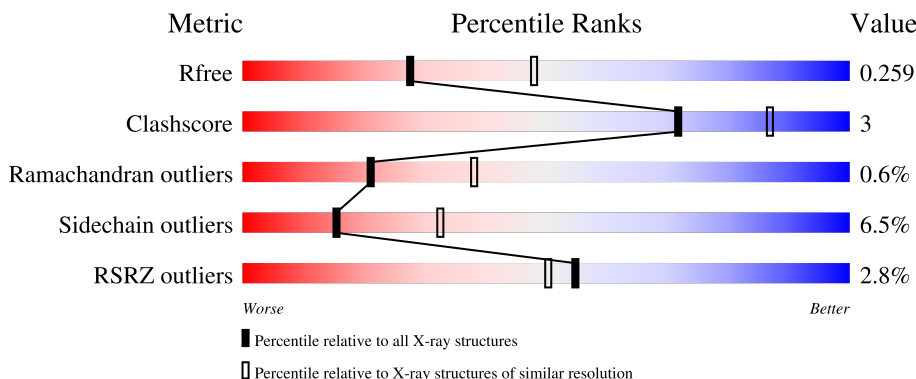
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	136	0.2% 69% 28%
1	E	136	0.2% 66% 6% 26%
1	K	136	2% 70% 27%
1	O	136	0.2% 68% 26%
2	B	103	3% 74% 7% 18%

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Mol	Chain	Length	Quality of chain
2	F	103	 6% 78% 5% 16%
2	L	103	 3% 68% 12% 17%
2	P	103	 13% 83% 6% 9%
3	C	130	 2% 78% 8% 14%
3	G	130	 2% 77% 18% 1%
3	M	130	 2% 74% 11% 15%
3	Q	130	 2% 72% 8% 18%
4	D	126	 2% 68% 7% 24%
4	H	126	 2% 64% 10% 24%
4	N	126	 67% 8% 23%
4	R	126	 2% 59% 17% 24%
5	I	338	 81% 17% 2%
6	J	338	 75% 24% 1%
7	S	213	 7% 28% 9% 61%

2 Entry composition

There are 11 unique types of molecules in this entry. The entry contains 27327 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 Histone H3.1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	98	Total	C	N	O	S	0	0	0
			807	508	156	139	4			
1	E	100	Total	C	N	O	S	0	0	0
			825	520	160	141	4			
1	K	99	Total	C	N	O	S	0	0	0
			816	514	158	140	4			
1	O	100	Total	C	N	O	S	0	0	0
			825	520	160	141	4			

- Molecule 2 is a protein called Histone H4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	84	Total	C	N	O	S	0	0	0
			673	424	133	115	1			
2	F	87	Total	C	N	O	S	0	0	0
			703	442	142	118	1			
2	L	85	Total	C	N	O	S	0	0	0
			683	430	136	116	1			
2	P	94	Total	C	N	O	S	0	0	0
			741	465	150	125	1			

- Molecule 3 is a protein called Histone H2A type 1-B/E.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	112	Total	C	N	O	0	0	0
			865	544	170	151			
3	G	107	Total	C	N	O	0	0	0
			828	523	162	143			
3	M	111	Total	C	N	O	0	0	0
			860	541	169	150			
3	Q	106	Total	C	N	O	0	0	0
			819	517	160	142			

- Molecule 4 is a protein called Histone H2B type 1-J.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	96	Total	C	N	O	S	0	0	0
			755	474	138	141	2			
4	H	96	Total	C	N	O	S	0	0	0
			755	474	138	141	2			
4	N	97	Total	C	N	O	S	0	0	0
			766	480	142	142	2			
4	R	96	Total	C	N	O	S	0	0	0
			755	474	138	141	2			

- Molecule 5 is a DNA chain called DNA (338-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	I	338	Total	C	N	O	P	0	0	0
			6923	3280	1316	1989	338			

- Molecule 6 is a DNA chain called DNA (338-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	J	338	Total	C	N	O	P	0	0	0
			6937	3298	1232	2069	338			

- Molecule 7 is a protein called Histone H1x.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
7	S	84	Total	C	N	O	0	0	0
			673	425	126	122			

- Molecule 8 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	1	Total	Cl	0	0
			1	1		
8	C	1	Total	Cl	0	0
			1	1		
8	E	1	Total	Cl	0	0
			1	1		
8	G	1	Total	Cl	0	0
			1	1		
8	K	1	Total	Cl	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	M	1	Total 1	Cl 1	0	0
8	O	1	Total 1	Cl 1	0	0
8	Q	1	Total 1	Cl 1	0	0

- Molecule 9 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	C	3	Total 3	Ca 3	0	0
9	D	1	Total 1	Ca 1	0	0
9	G	3	Total 3	Ca 3	0	0
9	I	28	Total 28	Ca 28	0	0
9	J	23	Total 23	Ca 23	0	0
9	M	2	Total 2	Ca 2	0	0
9	N	2	Total 2	Ca 2	0	0
9	Q	2	Total 2	Ca 2	0	0
9	R	1	Total 1	Ca 1	0	0

- Molecule 10 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	C	1	Total 1	K 1	0	0
10	I	6	Total 6	K 6	0	0
10	J	2	Total 2	K 2	0	0
10	M	1	Total 1	K 1	0	0

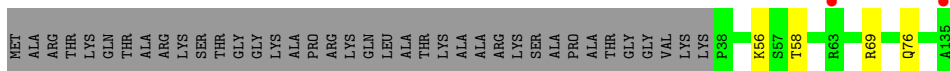
- Molecule 11 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
11	A	3	Total O 3 3	0	0
11	B	3	Total O 3 3	0	0
11	C	10	Total O 10 10	0	0
11	D	6	Total O 6 6	0	0
11	E	8	Total O 8 8	0	0
11	F	9	Total O 9 9	0	0
11	G	4	Total O 4 4	0	0
11	H	3	Total O 3 3	0	0
11	I	45	Total O 45 45	0	0
11	J	46	Total O 46 46	0	0
11	K	12	Total O 12 12	0	0
11	L	10	Total O 10 10	0	0
11	M	13	Total O 13 13	0	0
11	N	6	Total O 6 6	0	0
11	O	19	Total O 19 19	0	0
11	P	18	Total O 18 18	0	0
11	Q	12	Total O 12 12	0	0
11	R	8	Total O 8 8	0	0

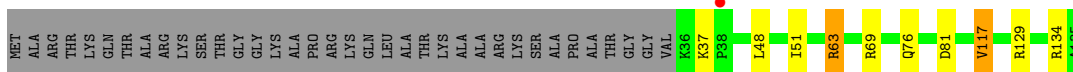
3 Residue-property plots [i](#)

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.

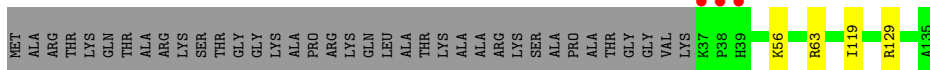
- Molecule 1: Histone H3.1



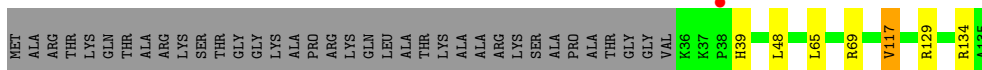
- Molecule 1: Histone H3.1



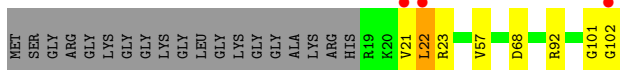
- Molecule 1: Histone H3.1



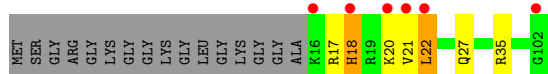
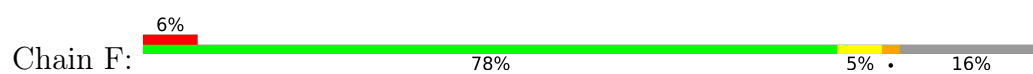
- Molecule 1: Histone H3.1



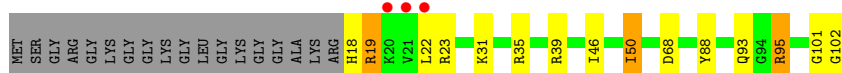
- Molecule 2: Histone H4



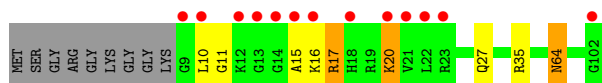
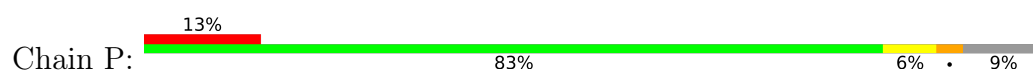
- Molecule 2: Histone H4



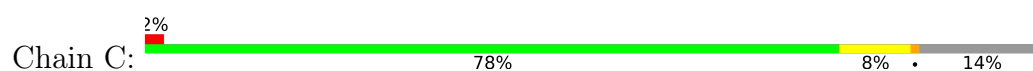
- Molecule 2: Histone H4



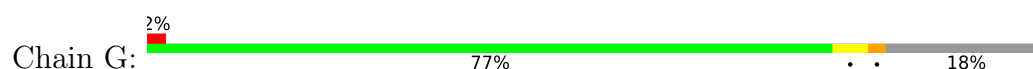
- Molecule 2: Histone H4



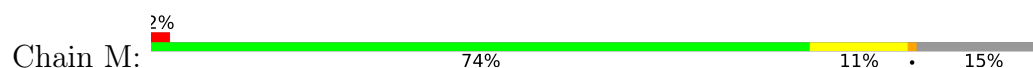
- Molecule 3: Histone H2A type 1-B/E



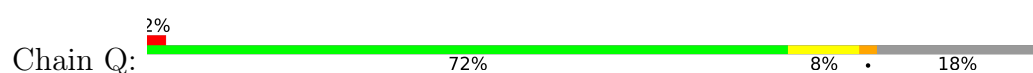
- Molecule 3: Histone H2A type 1-B/E



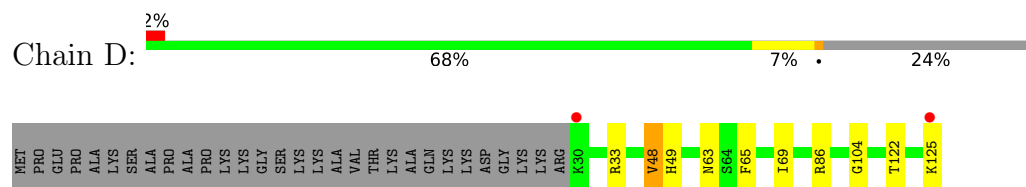
- Molecule 3: Histone H2A type 1-B/E



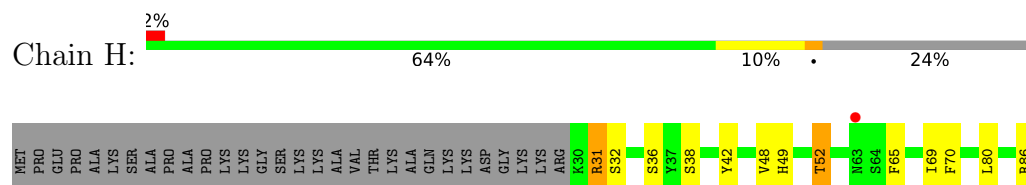
- Molecule 3: Histone H2A type 1-B/E



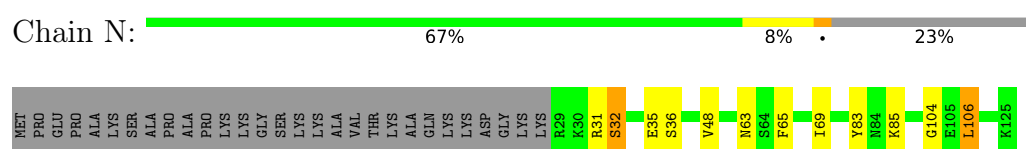
- Molecule 4: Histone H2B type 1-J



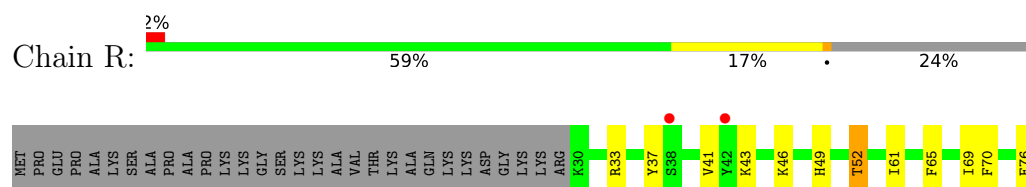
- Molecule 4: Histone H2B type 1-J



- Molecule 4: Histone H2B type 1-J



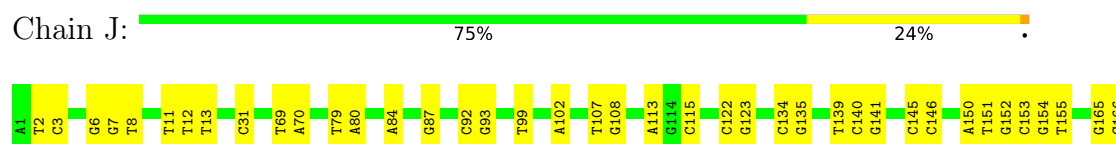
- Molecule 4: Histone H2B type 1-J

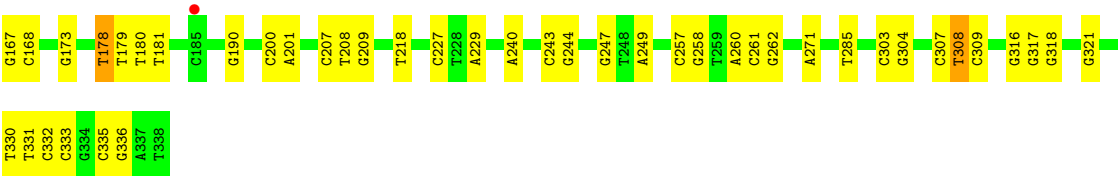


- Molecule 5: DNA (338-MER)

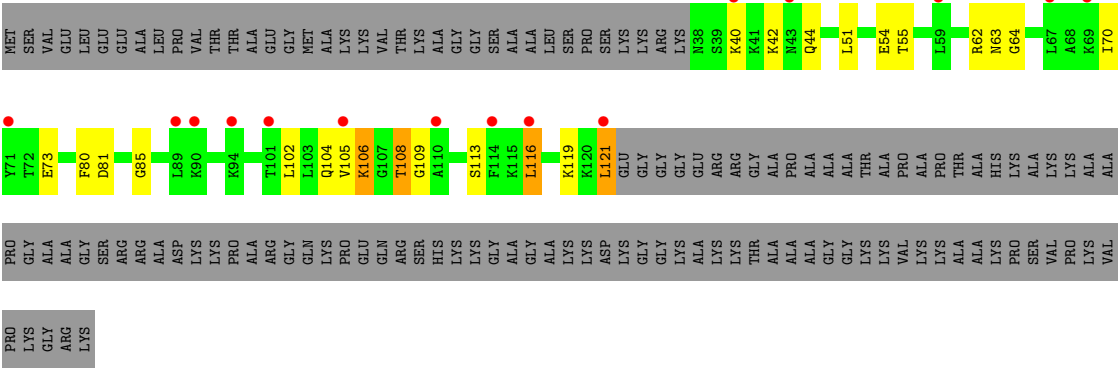


- Molecule 6: DNA (338-MER)





● Molecule 7: Histone H1x



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	103.50Å 101.41Å 215.67Å 90.00° 97.48° 90.00°	Depositor
Resolution (Å)	213.83 – 2.50 213.83 – 2.50	Depositor EDS
% Data completeness (in resolution range)	98.0 (213.83-2.50) 98.0 (213.83-2.50)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.36 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.8.0189	Depositor
R, R_{free}	0.204 , 0.256 0.209 , 0.259	Depositor DCC
R_{free} test set	2948 reflections (1.92%)	wwPDB-VP
Wilson B-factor (Å ²)	58.2	Xtriage
Anisotropy	0.029	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 43.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	27327	wwPDB-VP
Average B, all atoms (Å ²)	84.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.63% 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: CA, CL, K

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	0.74	0/819	0.97	0/1097
1	E	0.80	0/837	1.03	1/1120 (0.1%)
1	K	0.93	0/828	0.97	0/1109
1	O	0.96	0/837	1.01	1/1120 (0.1%)
2	B	0.85	0/680	0.96	1/908 (0.1%)
2	F	0.87	0/711	1.04	2/948 (0.2%)
2	L	0.99	1/691 (0.1%)	1.05	1/923 (0.1%)
2	P	1.04	0/749	1.10	1/997 (0.1%)
3	C	0.81	0/875	0.98	0/1179
3	G	0.80	0/838	0.95	1/1129 (0.1%)
3	M	0.91	0/870	1.03	0/1172
3	Q	0.84	0/829	0.98	1/1118 (0.1%)
4	D	0.85	0/766	1.01	0/1026
4	H	0.80	0/766	1.04	1/1026 (0.1%)
4	N	0.88	0/777	1.10	0/1040
4	R	0.82	0/766	1.10	0/1026
5	I	0.42	0/7778	0.90	20/11992 (0.2%)
6	J	0.43	0/7770	0.94	20/11998 (0.2%)
7	S	0.84	0/682	1.02	0/911
All	All	0.67	1/28869 (0.0%)	0.96	50/41839 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	L	31	LYS	C-O	-5.47	1.19	1.24

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	J	150	DA	C2'-C3'-O3'	8.15	123.73	111.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	I	74	DG	C2'-C3'-O3'	7.43	122.65	111.50
6	J	31	DC	C4'-C3'-O3'	-7.23	99.16	110.00
5	I	139	DG	C4'-C3'-O3'	7.14	120.71	110.00
5	I	188	DA	C2'-C3'-O3'	7.06	122.09	111.50
6	J	207	DC	C4'-C3'-O3'	-6.97	99.55	110.00
6	J	11	DT	C2'-C3'-O3'	6.96	121.94	111.50
5	I	332	DC	C2'-C3'-O3'	6.91	121.86	111.50
5	I	168	DG	N9-C1'-C2'	6.78	123.67	113.50
5	I	331	DA	C2'-C3'-O3'	6.64	121.47	111.50
5	I	171	DG	C4'-C3'-O3'	6.53	119.79	110.00
6	J	218	DT	C2'-C3'-O3'	-6.47	101.80	111.50
5	I	300	DA	C2'-C3'-O3'	-6.43	101.85	111.50
6	J	200	DC	C4'-C3'-O3'	-6.35	100.47	110.00
2	P	17	ARG	N-CA-C	6.35	119.92	110.52
5	I	51	DG	C2'-C3'-O3'	-6.31	102.03	111.50
6	J	113	DA	C4'-C3'-O3'	-6.31	100.54	110.00
1	O	117	VAL	N-CA-CB	-6.06	105.12	112.33
2	L	46	ILE	N-CA-C	6.00	117.33	107.24
6	J	308	DT	C4'-C3'-O3'	5.99	118.98	110.00
5	I	322	DT	C2'-C3'-O3'	-5.95	102.58	111.50
3	G	113	ALA	N-CA-C	5.89	118.18	111.11
6	J	99	DT	C2'-C3'-O3'	5.85	120.28	111.50
5	I	89	DC	C2'-C3'-O3'	-5.84	102.74	111.50
5	I	270	DA	C2'-C3'-O3'	-5.83	102.75	111.50
6	J	331	DT	C2'-C3'-O3'	5.81	120.22	111.50
4	H	38	SER	N-CA-C	5.76	118.30	111.33
6	J	318	DG	C2'-C3'-O3'	5.68	120.03	111.50
6	J	141	DG	C2'-C3'-O3'	-5.60	103.10	111.50
5	I	165	DG	C2'-C3'-O3'	5.57	119.86	111.50
6	J	307	DC	C2'-C3'-O3'	5.50	119.75	111.50
5	I	240	DA	C2'-C3'-O3'	5.49	119.74	111.50
5	I	240	DA	C4'-C3'-O3'	-5.46	101.81	110.00
6	J	249	DA	C2'-C3'-O3'	-5.45	103.33	111.50
3	Q	16	THR	N-CA-C	-5.45	100.43	108.99
6	J	260	DA	C4'-C3'-O3'	-5.44	101.85	110.00
6	J	178	DT	C2'-C3'-O3'	5.42	119.63	111.50
5	I	80	DA	C2'-C3'-O3'	-5.42	103.37	111.50
2	F	17	ARG	CA-C-N	5.29	131.65	121.54
2	F	17	ARG	C-N-CA	5.29	131.65	121.54
6	J	229	DA	C2'-C3'-O3'	-5.25	103.62	111.50
1	E	51	ILE	CB-CA-C	-5.24	105.27	111.97
5	I	31	DA	C4'-C3'-O3'	-5.23	102.15	110.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	I	201	DG	C2'-C3'-O3'	5.21	119.31	111.50
6	J	330	DT	C4'-C3'-O3'	-5.17	102.25	110.00
5	I	131	DA	C2'-C3'-O3'	-5.15	103.77	111.50
5	I	272	DC	C2'-C3'-O3'	5.02	119.03	111.50
6	J	227	DC	C2'-C3'-O3'	-5.02	103.97	111.50
6	J	87	DG	C2'-C3'-O3'	5.02	119.03	111.50
2	B	22	LEU	N-CA-C	5.00	115.90	108.60

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	807	0	844	5	0
1	E	825	0	869	5	0
1	K	816	0	856	2	0
1	O	825	0	869	2	0
2	B	673	0	722	3	0
2	F	703	0	755	3	0
2	L	683	0	729	9	0
2	P	741	0	796	5	0
3	C	865	0	928	6	0
3	G	828	0	892	3	0
3	M	860	0	923	7	0
3	Q	819	0	879	4	0
4	D	755	0	784	6	0
4	H	755	0	784	10	0
4	N	766	0	797	7	0
4	R	755	0	784	11	0
5	I	6923	0	3777	40	0
6	J	6937	0	3817	49	0
7	S	673	0	710	7	0
8	A	1	0	0	0	0
8	C	1	0	0	0	0
8	E	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	G	1	0	0	0	0
8	K	1	0	0	0	0
8	M	1	0	0	0	0
8	O	1	0	0	0	0
8	Q	1	0	0	0	0
9	C	3	0	0	0	0
9	D	1	0	0	0	0
9	G	3	0	0	0	0
9	I	28	0	0	0	0
9	J	23	0	0	0	0
9	M	2	0	0	0	0
9	N	2	0	0	0	0
9	Q	2	0	0	0	0
9	R	1	0	0	0	0
10	C	1	0	0	0	0
10	I	6	0	0	0	0
10	J	2	0	0	0	0
10	M	1	0	0	0	0
11	A	3	0	0	0	0
11	B	3	0	0	0	0
11	C	10	0	0	0	0
11	D	6	0	0	0	0
11	E	8	0	0	0	0
11	F	9	0	0	0	0
11	G	4	0	0	0	0
11	H	3	0	0	1	0
11	I	45	0	0	0	0
11	J	46	0	0	0	0
11	K	12	0	0	0	0
11	L	10	0	0	1	0
11	M	13	0	0	0	0
11	N	6	0	0	0	0
11	O	19	0	0	1	0
11	P	18	0	0	1	0
11	Q	12	0	0	0	0
11	R	8	0	0	0	0
All	All	27327	0	21515	149	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 3.

All (149) 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)
5:I:165:DG:H2''	5:I:166:DC:H5'	1.29	1.13
4:R:49:HIS:HB3	4:R:52:THR:HG23	1.56	0.85
5:I:165:DG:H2''	5:I:166:DC:C5'	2.08	0.82
5:I:335:DC:H1'	5:I:336:DG:H5''	1.65	0.77
2:L:95:ARG:NH1	11:L:201:HOH:O	2.10	0.75
5:I:335:DC:H1'	5:I:336:DG:C5'	2.18	0.73
4:H:49:HIS:HB3	4:H:52:THR:HG23	1.69	0.72
6:J:6:DG:H2''	6:J:7:DG:H5'	1.72	0.70
4:H:49:HIS:HB3	4:H:52:THR:CG2	2.21	0.70
1:E:76:GLN:HE22	2:F:22:LEU:HB3	1.59	0.67
4:D:122:THR:O	4:D:125:LYS:HD3	1.97	0.65
1:E:81:ASP:OD1	2:F:18:HIS:NE2	2.30	0.65
4:H:65:PHE:CE1	4:H:69:ILE:CD1	2.80	0.64
6:J:6:DG:H2''	6:J:7:DG:C5'	2.28	0.64
6:J:6:DG:H2''	6:J:7:DG:O4'	1.98	0.63
6:J:208:DT:H2''	6:J:209:DG:C8	2.33	0.63
4:N:65:PHE:CE1	4:N:69:ILE:CD1	2.82	0.63
6:J:180:DT:H4'	6:J:181:DT:OP1	1.98	0.62
4:D:65:PHE:CE1	4:D:69:ILE:CD1	2.84	0.60
5:I:171:DG:H4'	5:I:172:DC:OP1	2.01	0.60
4:R:65:PHE:CE1	4:R:69:ILE:CD1	2.86	0.59
1:K:119:ILE:HD12	2:L:50:ILE:HD13	1.85	0.59
5:I:232:DA:OP1	2:P:20:LYS:HB3	2.03	0.58
1:A:56:LYS:NZ	6:J:190:DG:OP1	2.36	0.58
7:S:108:THR:OG1	7:S:109:GLY:N	2.36	0.58
6:J:92:DC:H2''	6:J:93:DG:C8	2.39	0.57
7:S:55:THR:HG22	7:S:70:ILE:HG22	1.86	0.56
1:O:39:HIS:NE2	11:O:301:HOH:O	2.33	0.56
6:J:69:DT:H2''	6:J:70:DA:C8	2.40	0.55
5:I:167:DA:C2	6:J:173:DG:C2	2.94	0.55
3:M:120:THR:HG22	3:M:121:GLU:HG2	1.89	0.55
6:J:151:DT:H2''	6:J:152:DG:C8	2.43	0.54
3:C:115:LEU:HD23	1:E:117:VAL:HG22	1.90	0.54
5:I:190:DC:H2'	5:I:191:DC:C6	2.43	0.54
6:J:308:DT:H2''	6:J:309:DC:C6	2.43	0.54
6:J:2:DT:H5'	7:S:119:LYS:HE3	1.89	0.53
5:I:170:DA:H1'	5:I:171:DG:O4'	2.09	0.53
6:J:165:DG:H2''	6:J:166:DC:H5''	1.90	0.53
4:N:65:PHE:CE1	4:N:69:ILE:HD11	2.44	0.52
2:P:64:ASN:HD22	2:P:64:ASN:N	2.06	0.52
6:J:79:DT:H2''	6:J:80:DA:C8	2.45	0.52
4:R:49:HIS:HB3	4:R:52:THR:CG2	2.34	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:J:115:DC:OP1	4:N:32:SER:OG	2.21	0.52
4:H:31:ARG:NH1	6:J:285:DT:OP1	2.43	0.51
5:I:308:DG:H4'	5:I:309:DT:OP1	2.11	0.51
5:I:165:DG:C2'	5:I:166:DC:H5'	2.21	0.51
5:I:207:DT:H2''	5:I:208:DC:C6	2.46	0.50
6:J:303:DC:H2'	6:J:304:DG:C8	2.47	0.50
3:M:79:ILE:HG12	3:M:82:HIS:CE1	2.47	0.49
2:P:11:GLY:O	2:P:15:ALA:HB2	2.11	0.49
1:E:69:ARG:NH2	5:I:102:DA:OP2	2.46	0.49
5:I:23:DC:H2'	5:I:24:DG:C8	2.47	0.49
6:J:335:DC:H2'	6:J:336:DG:O4'	2.12	0.49
6:J:2:DT:H4'	6:J:3:DC:OP1	2.13	0.49
3:C:77:ARG:CZ	5:I:31:DA:H4'	2.43	0.49
3:C:81:ARG:NH2	3:C:107:VAL:O	2.35	0.49
5:I:85:DC:H1'	5:I:86:DT:H5'	1.95	0.49
3:Q:24:GLN:HG3	4:R:43:LYS:HB3	1.94	0.48
6:J:257:DC:H2''	6:J:258:DG:C8	2.48	0.48
1:A:76:GLN:HE21	2:B:22:LEU:CD1	2.26	0.48
3:M:81:ARG:NH2	3:M:107:VAL:O	2.39	0.48
5:I:81:DC:H2''	5:I:82:DG:C8	2.48	0.48
5:I:233:DC:H2''	5:I:234:DC:H5'	1.95	0.48
6:J:7:DG:H2''	6:J:8:DT:H72	1.96	0.48
6:J:102:DA:OP2	1:O:69:ARG:NH2	2.46	0.48
4:H:49:HIS:CB	4:H:52:THR:HG23	2.42	0.48
5:I:122:DC:H2'	5:I:123:DT:H71	1.95	0.48
3:M:31:HIS:CD2	3:M:48:PRO:HG3	2.49	0.47
3:C:77:ARG:NE	5:I:31:DA:H4'	2.28	0.47
1:A:69:ARG:HH22	6:J:271:DA:P	2.37	0.47
6:J:316:DG:H1'	6:J:317:DG:H5'	1.96	0.47
5:I:12:DA:C2	5:I:13:DA:C2	3.03	0.47
6:J:122:DC:H2''	6:J:123:DG:C8	2.50	0.47
6:J:178:DT:H1'	6:J:179:DT:O4'	2.15	0.47
5:I:152:DA:H2'	5:I:153:DT:C6	2.49	0.47
2:L:19:ARG:HG2	2:L:23:ARG:HB2	1.97	0.47
3:M:32:ARG:NH1	4:N:35:GLU:OE2	2.44	0.47
3:Q:102:ILE:HG23	4:R:61:ILE:HD13	1.97	0.46
4:D:33:ARG:HH12	5:I:40:DA:H5'	1.81	0.46
5:I:187:DC:H2''	5:I:188:DA:O5'	2.15	0.46
2:B:68:ASP:OD2	2:B:92:ARG:NH1	2.47	0.46
6:J:243:DC:H2''	6:J:244:DG:C8	2.50	0.46
2:L:101:GLY:O	2:L:102:GLY:C	2.59	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:35:ARG:NH2	5:I:93:DC:OP2	2.46	0.46
5:I:149:DC:H2'	5:I:150:DA:C8	2.51	0.45
3:G:84:GLN:NE2	3:G:106:GLY:O	2.45	0.45
5:I:186:DG:C2	6:J:154:DG:C2	3.05	0.45
4:H:42:TYR:OH	6:J:201:DA:OP2	2.27	0.45
6:J:179:DT:H1'	6:J:180:DT:H5'	1.97	0.45
4:N:65:PHE:CZ	4:N:69:ILE:CD1	3.00	0.45
5:I:189:DT:OP1	1:K:56:LYS:NZ	2.39	0.45
6:J:12:DT:H1'	6:J:13:DT:H5'	1.99	0.45
4:R:70:PHE:CD1	4:R:70:PHE:C	2.94	0.45
6:J:139:DT:H2''	6:J:140:DC:C6	2.52	0.45
4:H:52:THR:HG22	11:H:202:HOH:O	2.15	0.45
5:I:157:DA:H2'	5:I:158:DA:C8	2.52	0.45
7:S:106:LYS:HD2	7:S:113:SER:OG	2.17	0.44
6:J:154:DG:C2	6:J:155:DT:C2	3.05	0.44
5:I:19:DA:C2	6:J:321:DG:C2	3.06	0.44
4:R:78:SER:HA	4:R:89:ILE:HD11	1.99	0.44
6:J:153:DC:H2''	6:J:154:DG:C5'	2.47	0.44
3:Q:81:ARG:NH2	3:Q:107:VAL:O	2.45	0.44
3:G:73:ASN:N	3:G:73:ASN:HD22	2.16	0.44
5:I:335:DC:OP2	5:I:335:DC:H2'	2.18	0.44
7:S:80:PHE:CE1	7:S:85:GLY:HA3	2.53	0.44
5:I:187:DC:C6	5:I:187:DC:H5'	2.53	0.43
3:M:92:GLU:HB3	4:N:106:LEU:HD22	2.00	0.43
2:P:35:ARG:HD2	11:P:217:HOH:O	2.16	0.43
3:C:79:ILE:HG12	3:C:82:HIS:CE1	2.54	0.43
6:J:107:DT:H2''	6:J:108:DG:H5'	2.01	0.43
6:J:139:DT:H2''	6:J:140:DC:O5'	2.18	0.43
1:A:58:THR:HG21	3:G:81:ARG:HB2	2.00	0.42
2:L:35:ARG:O	2:L:39:ARG:HG2	2.19	0.42
2:L:18:HIS:CG	2:L:19:ARG:H	2.37	0.42
6:J:261:DC:H2''	6:J:262:DG:C8	2.55	0.42
4:R:76:GLU:OE2	4:R:79:ARG:NH1	2.52	0.42
2:B:101:GLY:O	2:B:102:GLY:C	2.63	0.42
1:A:69:ARG:NH2	6:J:271:DA:OP2	2.53	0.42
5:I:335:DC:H1'	5:I:336:DG:H5'	1.97	0.42
4:R:102:LEU:HB2	4:R:107:ALA:HB2	2.01	0.42
6:J:6:DG:C2'	6:J:7:DG:O4'	2.67	0.42
4:D:48:VAL:HG12	4:D:49:HIS:CD2	2.54	0.42
5:I:265:DG:H2''	5:I:266:DT:OP2	2.20	0.42
4:H:70:PHE:CD1	4:H:70:PHE:C	2.98	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:63:ARG:CZ	6:J:240:DA:H4'	2.50	0.41
5:I:186:DG:N2	6:J:154:DG:C2	2.88	0.41
6:J:145:DC:H2''	6:J:146:DC:OP2	2.19	0.41
5:I:180:DA:H5'	5:I:180:DA:C8	2.55	0.41
6:J:153:DC:H2''	6:J:154:DG:O5'	2.20	0.41
2:L:19:ARG:CD	2:L:23:ARG:HB2	2.51	0.41
4:H:65:PHE:CZ	4:H:69:ILE:HD13	2.55	0.41
4:H:65:PHE:CE1	4:H:69:ILE:HD11	2.56	0.41
3:M:88:ARG:HD3	3:M:88:ARG:HA	1.88	0.41
6:J:167:DG:H1'	6:J:168:DC:H5'	2.02	0.41
6:J:332:DC:H2''	6:J:333:DC:H5''	2.03	0.41
7:S:63:ASN:OD1	7:S:106:LYS:NZ	2.54	0.41
5:I:3:DC:H5'	5:I:3:DC:H6	1.85	0.41
5:I:91:DA:N6	6:J:247:DG:C6	2.89	0.41
2:L:88:TYR:CE1	4:N:83:TYR:CE1	3.09	0.41
4:R:37:TYR:O	4:R:41:VAL:HG23	2.21	0.41
5:I:335:DC:H2''	5:I:336:DG:H5'	2.03	0.40
3:C:62:ILE:HD12	4:D:65:PHE:CZ	2.56	0.40
2:L:68:ASP:OD2	2:L:93:GLN:NE2	2.55	0.40
7:S:121:LEU:HD12	7:S:121:LEU:H	1.87	0.40
4:D:65:PHE:CE1	4:D:69:ILE:HD11	2.55	0.40
6:J:134:DC:H2'	6:J:135:DG:C8	2.56	0.40
2:P:11:GLY:C	2:P:15:ALA:HB2	2.46	0.40
5:I:256:DG:C2	6:J:84:DA:C2	3.09	0.40
3:Q:67:GLY:HA3	4:R:49:HIS:CD2	2.57	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	96/136 (71%)	96 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	98/136 (72%)	97 (99%)	1 (1%)	0	100	100
1	K	97/136 (71%)	97 (100%)	0	0	100	100
1	O	98/136 (72%)	97 (99%)	1 (1%)	0	100	100
2	B	82/103 (80%)	79 (96%)	2 (2%)	1 (1%)	10	20
2	F	85/103 (82%)	79 (93%)	4 (5%)	2 (2%)	4	8
2	L	83/103 (81%)	80 (96%)	3 (4%)	0	100	100
2	P	92/103 (89%)	86 (94%)	6 (6%)	0	100	100
3	C	110/130 (85%)	106 (96%)	4 (4%)	0	100	100
3	G	105/130 (81%)	100 (95%)	4 (4%)	1 (1%)	12	24
3	M	109/130 (84%)	103 (94%)	5 (5%)	1 (1%)	14	27
3	Q	104/130 (80%)	102 (98%)	2 (2%)	0	100	100
4	D	94/126 (75%)	93 (99%)	0	1 (1%)	11	22
4	H	94/126 (75%)	90 (96%)	4 (4%)	0	100	100
4	N	95/126 (75%)	90 (95%)	4 (4%)	1 (1%)	11	22
4	R	94/126 (75%)	91 (97%)	2 (2%)	1 (1%)	11	22
7	S	82/213 (38%)	72 (88%)	8 (10%)	2 (2%)	4	8
All	All	1618/2193 (74%)	1558 (96%)	50 (3%)	10 (1%)	21	38

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	23	ARG
2	F	18	HIS
4	R	33	ARG
4	D	104	GLY
3	M	119	LYS
4	N	104	GLY
7	S	64	GLY
2	F	20	LYS
3	G	118	LYS
7	S	116	LEU

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	85/111 (77%)	85 (100%)	0	100	100
1	E	87/111 (78%)	81 (93%)	6 (7%)	14	30
1	K	86/111 (78%)	84 (98%)	2 (2%)	44	72
1	O	87/111 (78%)	82 (94%)	5 (6%)	18	39
2	B	69/79 (87%)	67 (97%)	2 (3%)	37	65
2	F	72/79 (91%)	69 (96%)	3 (4%)	26	52
2	L	70/79 (89%)	66 (94%)	4 (6%)	18	39
2	P	74/79 (94%)	68 (92%)	6 (8%)	11	23
3	C	88/100 (88%)	83 (94%)	5 (6%)	18	39
3	G	85/100 (85%)	82 (96%)	3 (4%)	32	58
3	M	88/100 (88%)	84 (96%)	4 (4%)	24	49
3	Q	84/100 (84%)	76 (90%)	8 (10%)	8	17
4	D	82/105 (78%)	79 (96%)	3 (4%)	30	57
4	H	82/105 (78%)	73 (89%)	9 (11%)	6	13
4	N	83/105 (79%)	76 (92%)	7 (8%)	10	22
4	R	82/105 (78%)	75 (92%)	7 (8%)	10	22
7	S	73/157 (46%)	58 (80%)	15 (20%)	1	2
All	All	1377/1737 (79%)	1288 (94%)	89 (6%)	15	32

All (89) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	21	VAL
2	B	57	VAL
3	C	24	GLN
3	C	81	ARG
3	C	118	LYS
3	C	119	LYS
3	C	120	THR
4	D	48	VAL
4	D	63	ASN
4	D	86	ARG
1	E	37	LYS

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Mol	Chain	Res	Type
1	E	48	LEU
1	E	63	ARG
1	E	117	VAL
1	E	129	ARG
1	E	134	ARG
2	F	21	VAL
2	F	22	LEU
2	F	27	GLN
3	G	29	ARG
3	G	81	ARG
3	G	118	LYS
4	H	31	ARG
4	H	32	SER
4	H	36	SER
4	H	48	VAL
4	H	52	THR
4	H	80	LEU
4	H	86	ARG
4	H	106	LEU
4	H	125	LYS
1	K	63	ARG
1	K	129	ARG
2	L	19	ARG
2	L	22	LEU
2	L	50	ILE
2	L	95	ARG
3	M	24	GLN
3	M	73	ASN
3	M	74	LYS
3	M	81	ARG
4	N	31	ARG
4	N	32	SER
4	N	36	SER
4	N	48	VAL
4	N	63	ASN
4	N	85	LYS
4	N	106	LEU
1	O	48	LEU
1	O	65	LEU
1	O	117	VAL
1	O	129	ARG
1	O	134	ARG

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Mol	Chain	Res	Type
2	P	10	LEU
2	P	16	LYS
2	P	17	ARG
2	P	20	LYS
2	P	27	GLN
2	P	64	ASN
3	Q	15	LYS
3	Q	24	GLN
3	Q	29	ARG
3	Q	64	GLU
3	Q	71	ARG
3	Q	74	LYS
3	Q	76	THR
3	Q	81	ARG
4	R	46	LYS
4	R	52	THR
4	R	80	LEU
4	R	85	LYS
4	R	86	ARG
4	R	106	LEU
4	R	125	LYS
7	S	40	LYS
7	S	42	LYS
7	S	44	GLN
7	S	51	LEU
7	S	54	GLU
7	S	62	ARG
7	S	73	GLU
7	S	81	ASP
7	S	102	LEU
7	S	104	GLN
7	S	105	VAL
7	S	106	LYS
7	S	108	THR
7	S	116	LEU
7	S	121	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	76	GLN
1	A	85	GLN

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Mol	Chain	Res	Type
3	C	31	HIS
3	C	104	GLN
4	D	49	HIS
4	D	67	ASN
4	D	95	GLN
4	D	109	HIS
1	E	39	HIS
1	E	76	GLN
2	F	27	GLN
3	G	73	ASN
3	G	104	GLN
1	K	39	HIS
2	L	64	ASN
3	M	31	HIS
3	M	89	ASN
4	N	47	GLN
1	O	76	GLN
1	O	85	GLN
2	P	27	GLN
2	P	64	ASN
3	Q	24	GLN
4	R	47	GLN
7	S	38	ASN
7	S	44	GLN
7	S	98	GLN
7	S	99	ASN
7	S	104	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 83 ligands modelled in this entry, 83 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

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2	OWAB(Å ²)	Q < 0.9
1	A	98/136 (72%)	0.07	2 (2%) 65 61	45, 59, 90, 110	0
1	E	100/136 (73%)	-0.18	1 (1%) 79 76	39, 53, 88, 133	0
1	K	99/136 (72%)	-0.08	3 (3%) 52 48	33, 46, 86, 130	0
1	O	100/136 (73%)	-0.35	1 (1%) 79 76	32, 44, 76, 127	0
2	B	84/103 (81%)	0.06	3 (3%) 46 41	42, 55, 124, 161	0
2	F	87/103 (84%)	-0.10	6 (6%) 23 20	39, 50, 137, 162	0
2	L	85/103 (82%)	-0.12	3 (3%) 47 42	36, 46, 126, 174	0
2	P	94/103 (91%)	0.10	13 (13%) 6 5	34, 44, 157, 174	0
3	C	112/130 (86%)	-0.02	3 (2%) 56 51	41, 53, 115, 151	0
3	G	107/130 (82%)	-0.04	3 (2%) 55 50	44, 57, 96, 132	0
3	M	111/130 (85%)	-0.10	3 (2%) 56 51	36, 45, 97, 144	0
3	Q	106/130 (81%)	-0.18	2 (1%) 66 62	37, 50, 81, 128	0
4	D	96/126 (76%)	-0.00	2 (2%) 63 59	43, 55, 109, 129	0
4	H	96/126 (76%)	0.05	2 (2%) 63 59	45, 57, 99, 134	0
4	N	97/126 (76%)	-0.16	0 100 100	34, 47, 96, 144	0
4	R	96/126 (76%)	0.11	3 (3%) 51 47	38, 49, 93, 127	0
5	I	338/338 (100%)	-0.05	0 100 100	55, 94, 198, 230	0
6	J	338/338 (100%)	-0.02	1 (0%) 90 87	53, 96, 204, 251	0
7	S	84/213 (39%)	1.30	15 (17%) 3 3	112, 147, 171, 188	0
All	All	2328/2869 (81%)	-0.00	66 (2%) 55 50	32, 60, 154, 251	0

All (66) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	L	22	LEU	4.7

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Mol	Chain	Res	Type	RSRZ
2	P	21	VAL	4.6
2	P	10	LEU	4.4
1	K	38	PRO	4.0
2	L	21	VAL	3.8
3	Q	14	ALA	3.7
2	B	21	VAL	3.6
2	F	21	VAL	3.6
7	S	121	LEU	3.5
3	M	119	LYS	3.4
7	S	59	LEU	3.4
2	P	12	LYS	3.3
2	P	16	LYS	3.3
2	P	13	GLY	3.3
2	P	9	GLY	3.1
3	Q	119	LYS	3.0
2	F	16	LYS	3.0
2	P	102	GLY	2.9
2	B	22	LEU	2.9
2	P	14	GLY	2.9
2	P	15	ALA	2.9
3	G	14	ALA	2.9
7	S	116	LEU	2.9
4	D	30	LYS	2.8
2	P	22	LEU	2.8
3	G	13	LYS	2.7
2	P	20	LYS	2.6
3	M	120	THR	2.6
2	F	20	LYS	2.6
1	K	37	LYS	2.6
7	S	67	LEU	2.6
3	M	12	ALA	2.5
2	L	20	LYS	2.5
2	P	18	HIS	2.5
7	S	105	VAL	2.5
2	F	22	LEU	2.4
3	C	10	ALA	2.4
1	E	38	PRO	2.4
3	C	12	ALA	2.4
2	B	102	GLY	2.4
4	R	42	TYR	2.4
4	R	38	SER	2.3
7	S	94	LYS	2.3

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Mol	Chain	Res	Type	RSRZ
7	S	71	TYR	2.3
4	H	63	ASN	2.3
4	R	82	HIS	2.3
7	S	114	PHE	2.3
3	G	36	LYS	2.3
4	H	125	LYS	2.3
1	A	63	ARG	2.3
1	K	39	HIS	2.2
6	J	185	DC	2.2
1	A	135	ALA	2.2
7	S	43	ASN	2.2
3	C	119	LYS	2.2
1	O	38	PRO	2.2
7	S	69	LYS	2.2
2	F	18	HIS	2.1
7	S	89	LEU	2.1
7	S	110	ALA	2.1
2	F	102	GLY	2.1
4	D	125	LYS	2.1
7	S	40	LYS	2.1
2	P	23	ARG	2.0
7	S	90	LYS	2.0
7	S	101	THR	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no 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.

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
9	CA	J	418	1/1	0.80	0.10	101,101,101,101	0
9	CA	J	402	1/1	0.81	0.09	111,111,111,111	0
9	CA	G	202	1/1	0.82	0.12	117,117,117,117	0
9	CA	J	410	1/1	0.82	0.09	103,103,103,103	0
9	CA	I	401	1/1	0.82	0.09	110,110,110,110	0
10	K	I	433	1/1	0.82	0.21	114,114,114,114	0
9	CA	I	416	1/1	0.83	0.10	115,115,115,115	0
9	CA	G	203	1/1	0.84	0.13	102,102,102,102	0
9	CA	I	417	1/1	0.84	0.12	108,108,108,108	0
9	CA	I	424	1/1	0.86	0.14	100,100,100,100	0
9	CA	I	427	1/1	0.86	0.09	102,102,102,102	0
9	CA	I	420	1/1	0.87	0.08	108,108,108,108	0
9	CA	I	426	1/1	0.87	0.08	104,104,104,104	0
9	CA	J	408	1/1	0.87	0.08	104,104,104,104	0
10	K	I	434	1/1	0.87	0.13	110,110,110,110	0
9	CA	I	405	1/1	0.88	0.07	102,102,102,102	0
9	CA	I	415	1/1	0.88	0.08	102,102,102,102	0
9	CA	G	201	1/1	0.88	0.09	97,97,97,97	0
9	CA	J	421	1/1	0.88	0.09	100,100,100,100	0
9	CA	J	422	1/1	0.88	0.11	95,95,95,95	0
9	CA	R	201	1/1	0.88	0.08	97,97,97,97	0
9	CA	I	402	1/1	0.88	0.12	103,103,103,103	0
9	CA	I	418	1/1	0.88	0.09	97,97,97,97	0
9	CA	I	414	1/1	0.89	0.08	111,111,111,111	0
9	CA	I	422	1/1	0.89	0.08	91,91,91,91	0
9	CA	J	423	1/1	0.89	0.09	100,100,100,100	0
9	CA	J	413	1/1	0.89	0.07	106,106,106,106	0
9	CA	I	419	1/1	0.89	0.13	109,109,109,109	0
9	CA	J	420	1/1	0.89	0.08	111,111,111,111	0
9	CA	I	413	1/1	0.90	0.14	98,98,98,98	0
9	CA	I	423	1/1	0.90	0.08	95,95,95,95	0
9	CA	N	202	1/1	0.90	0.23	83,83,83,83	0
9	CA	Q	202	1/1	0.90	0.08	91,91,91,91	0
9	CA	I	412	1/1	0.90	0.07	92,92,92,92	0
9	CA	I	425	1/1	0.90	0.07	104,104,104,104	0
9	CA	J	409	1/1	0.90	0.09	102,102,102,102	0
10	K	C	205	1/1	0.91	0.08	93,93,93,93	0
10	K	I	432	1/1	0.91	0.07	99,99,99,99	0
9	CA	M	202	1/1	0.91	0.12	94,94,94,94	0
9	CA	J	419	1/1	0.91	0.07	101,101,101,101	0
9	CA	J	411	1/1	0.92	0.08	94,94,94,94	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
10	K	I	429	1/1	0.92	0.08	80,80,80,80	0
10	K	I	430	1/1	0.92	0.10	80,80,80,80	0
9	CA	I	421	1/1	0.92	0.12	102,102,102,102	0
9	CA	I	409	1/1	0.92	0.08	85,85,85,85	0
9	CA	J	404	1/1	0.92	0.10	93,93,93,93	0
9	CA	J	406	1/1	0.93	0.07	83,83,83,83	0
8	CL	K	201	1/1	0.93	0.14	64,64,64,64	0
9	CA	I	404	1/1	0.93	0.08	86,86,86,86	0
9	CA	J	416	1/1	0.93	0.08	95,95,95,95	0
10	K	M	204	1/1	0.93	0.07	80,80,80,80	0
9	CA	I	403	1/1	0.94	0.06	102,102,102,102	0
9	CA	N	201	1/1	0.94	0.06	84,84,84,84	0
9	CA	J	412	1/1	0.94	0.08	90,90,90,90	0
9	CA	I	428	1/1	0.94	0.24	82,82,82,82	0
9	CA	I	406	1/1	0.94	0.10	106,106,106,106	0
8	CL	E	201	1/1	0.94	0.14	66,66,66,66	0
9	CA	J	407	1/1	0.95	0.06	99,99,99,99	0
9	CA	C	203	1/1	0.95	0.16	87,87,87,87	0
9	CA	Q	201	1/1	0.95	0.08	98,98,98,98	0
9	CA	C	202	1/1	0.95	0.11	96,96,96,96	0
9	CA	J	405	1/1	0.95	0.06	112,112,112,112	0
10	K	J	424	1/1	0.95	0.10	77,77,77,77	0
9	CA	I	407	1/1	0.95	0.07	96,96,96,96	0
8	CL	A	201	1/1	0.96	0.07	72,72,72,72	0
9	CA	I	410	1/1	0.96	0.06	85,85,85,85	0
8	CL	Q	203	1/1	0.96	0.07	51,51,51,51	0
9	CA	C	201	1/1	0.96	0.07	87,87,87,87	0
9	CA	J	414	1/1	0.96	0.10	95,95,95,95	0
10	K	I	431	1/1	0.96	0.07	97,97,97,97	0
9	CA	M	201	1/1	0.96	0.11	91,91,91,91	0
9	CA	J	415	1/1	0.96	0.08	82,82,82,82	0
9	CA	I	408	1/1	0.96	0.05	91,91,91,91	0
9	CA	J	417	1/1	0.96	0.09	85,85,85,85	0
9	CA	J	403	1/1	0.96	0.06	79,79,79,79	0
9	CA	I	411	1/1	0.97	0.04	83,83,83,83	0
10	K	J	425	1/1	0.97	0.05	77,77,77,77	0
9	CA	J	401	1/1	0.97	0.08	88,88,88,88	0
9	CA	D	201	1/1	0.98	0.04	80,80,80,80	0
8	CL	G	204	1/1	0.98	0.04	61,61,61,61	0
8	CL	O	201	1/1	0.98	0.08	49,49,49,49	0
8	CL	M	203	1/1	0.99	0.13	48,48,48,48	0
8	CL	C	204	1/1	0.99	0.09	55,55,55,55	0

6.5 Other polymers [i](#)

There are no such residues in this entry.