



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 12, 2021 – 06:12 PM JST

PDB ID : 6L9Z  
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](mailto: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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

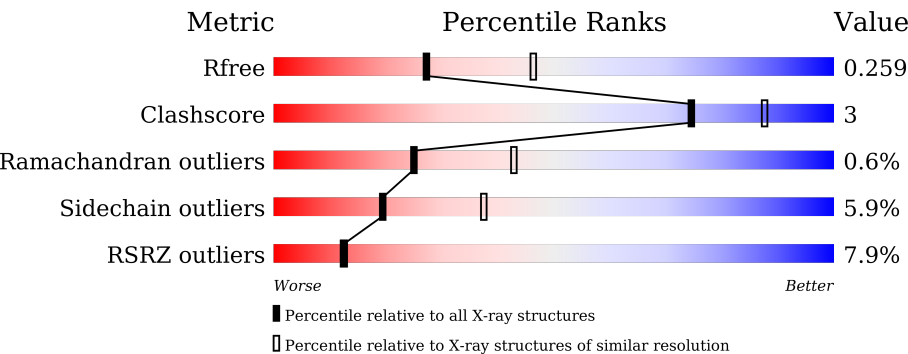
MolProbity : 4.02b-467  
Xtriage (Phenix) : 1.13  
EDS : 2.23.1  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.23.1

# 1 Overall quality at a glance i

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))
$R_{free}$	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (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  $\geq 3$ , 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	<div><div>2%</div><div><div></div><div>68%</div><div></div><div>28%</div></div><div></div></div>
1	E	136	<div><div></div><div><div></div><div>65%</div><div></div><div>26%</div></div><div><div></div><div>6%</div><div></div></div></div>
1	K	136	<div><div>3%</div><div><div></div><div>68%</div><div></div><div>27%</div></div><div></div></div>
1	O	136	<div><div>%</div><div><div></div><div>69%</div><div></div><div>26%</div></div><div></div></div>
2	B	103	<div><div>5%</div><div><div></div><div>77%</div><div></div><div>18%</div></div><div></div></div>
2	F	103	<div><div>5%</div><div><div></div><div>79%</div><div></div><div>16%</div></div><div></div></div>

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
2	L	103	
2	P	103	
3	C	130	
3	G	130	
3	M	130	
3	Q	130	
4	D	126	
4	H	126	
4	N	126	
4	R	126	
5	I	338	
6	J	338	
7	S	213	

## 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 (three-letter code: 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		

*Continued on next page...*

*Continued from previous page...*

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 (three-letter code: 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 (three-letter code: 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

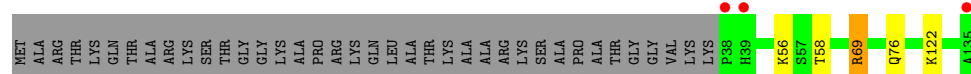
- Molecule 11 is water.

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

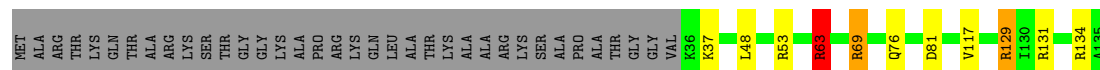
### 3 Residue-property plots [i](#)

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.

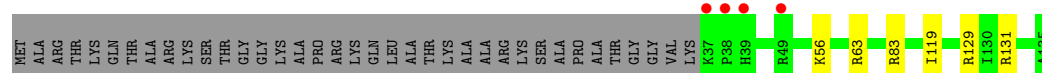
- Molecule 1: Histone H3.1



- Molecule 1: Histone H3.1



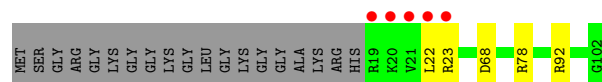
- Molecule 1: Histone H3.1



- Molecule 1: Histone H3.1

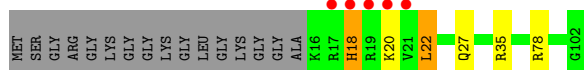
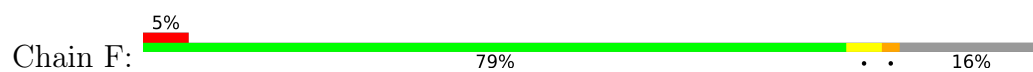


- Molecule 2: Histone H4

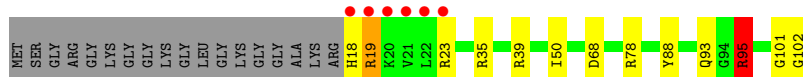


- Molecule 2: Histone H4

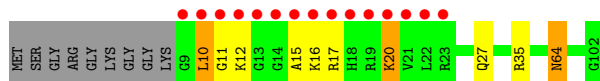
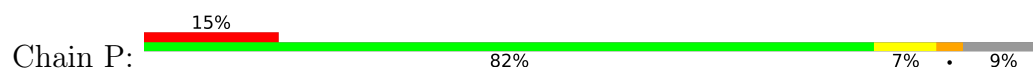




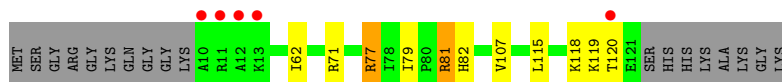
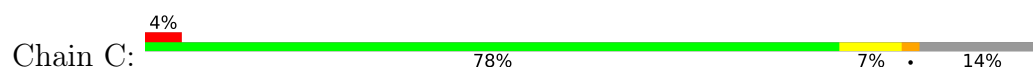
- Molecule 2: Histone H4



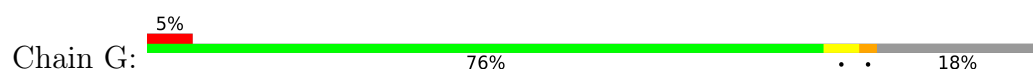
- Molecule 2: Histone H4



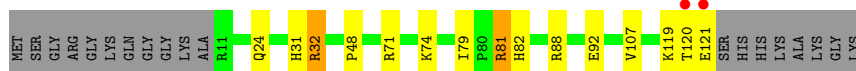
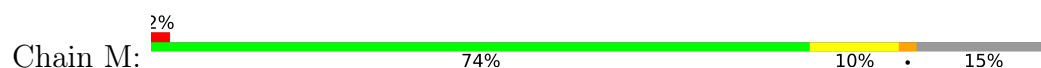
- Molecule 3: Histone H2A type 1-B/E



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



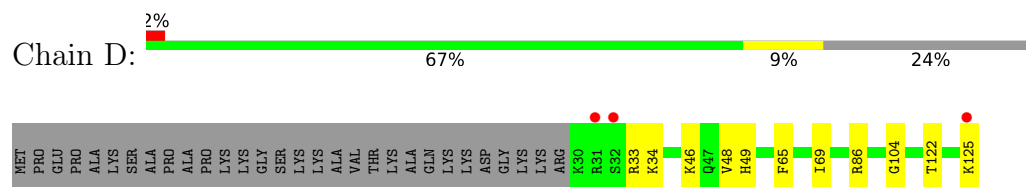
- Molecule 3: Histone H2A type 1-B/E



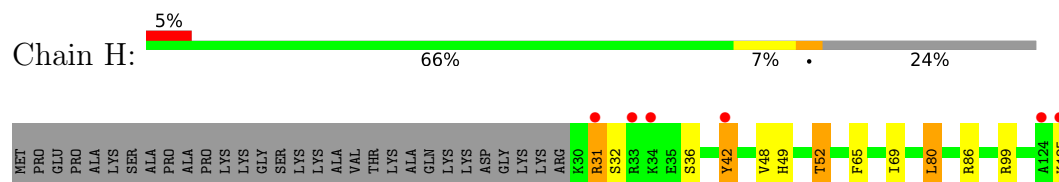
- Molecule 3: Histone H2A type 1-B/E



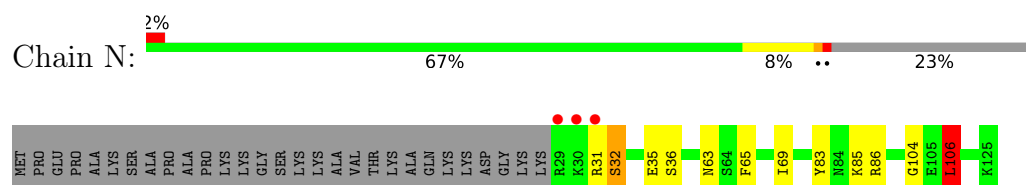
- Molecule 4: Histone H2B type 1-J



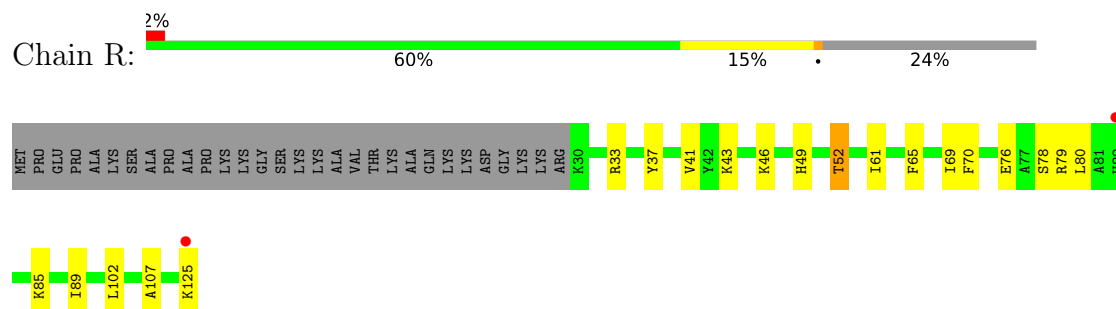
- Molecule 4: Histone H2B type 1-J



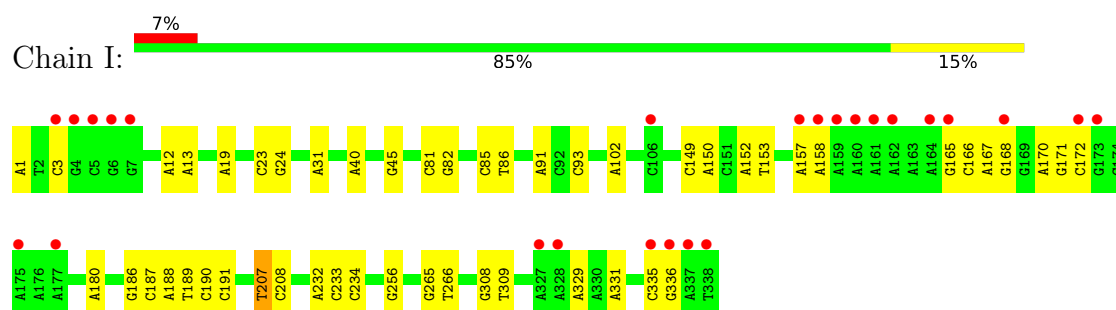
- Molecule 4: Histone H2B type 1-J



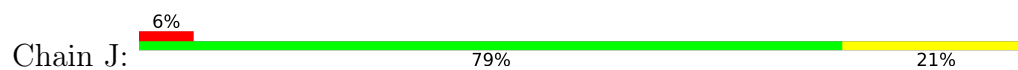
- Molecule 4: Histone H2B type 1-J

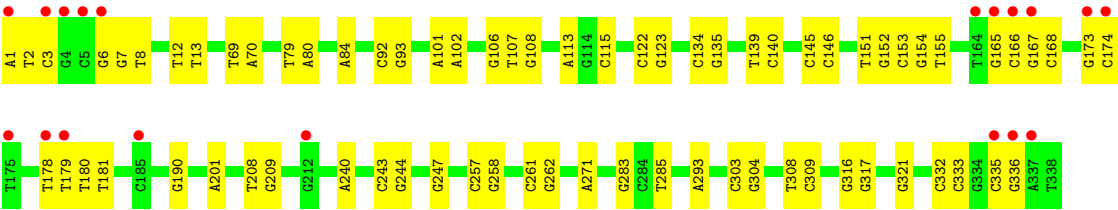


- Molecule 5: DNA (338-MER)

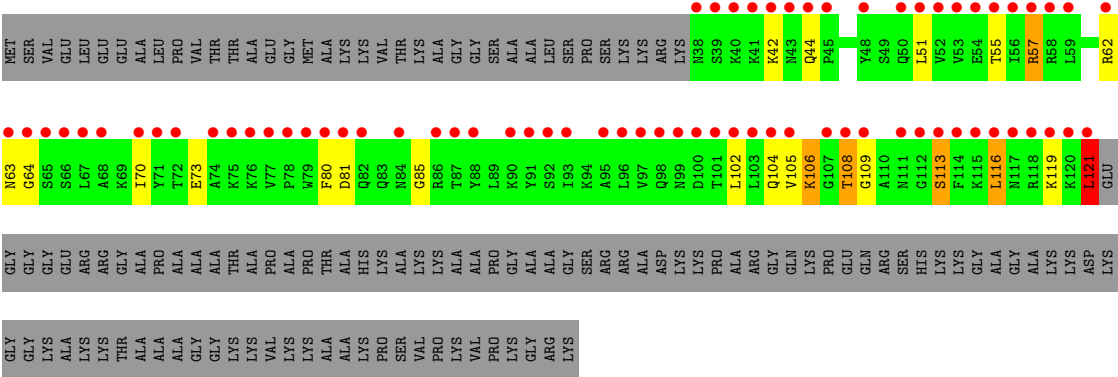


- Molecule 6: DNA (338-MER)





● Molecule 7: Histone H1x



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	103.50Å 101.41Å 215.67Å 90.00° 97.48° 90.00°	Depositor
Resolution (Å)	213.83 – 2.50 91.63 – 2.50	Depositor EDS
% Data completeness (in resolution range)	98.0 (213.83-2.50) 98.0 (91.63-2.50)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	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 (1.96%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	58.2	Xtriage
Anisotropy	0.029	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 48.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	27327	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	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.*

<sup>1</sup>Intensities estimated from amplitudes.

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

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CL, K, CA

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	0.53	0/819	0.86	1/1097 (0.1%)
1	E	0.60	0/837	0.97	5/1120 (0.4%)
1	K	0.66	0/828	0.94	2/1109 (0.2%)
1	O	0.67	0/837	0.91	3/1120 (0.3%)
2	B	0.59	0/680	0.89	1/908 (0.1%)
2	F	0.60	0/711	0.93	1/948 (0.1%)
2	L	0.67	0/691	0.97	2/923 (0.2%)
2	P	0.72	0/749	1.01	1/997 (0.1%)
3	C	0.56	0/875	0.82	3/1179 (0.3%)
3	G	0.56	0/838	0.77	0/1129
3	M	0.62	0/870	0.91	3/1172 (0.3%)
3	Q	0.59	0/829	0.88	3/1118 (0.3%)
4	D	0.64	0/766	0.80	0/1026
4	H	0.59	0/766	0.84	2/1026 (0.2%)
4	N	0.71	0/777	0.89	1/1040 (0.1%)
4	R	0.67	0/766	0.83	0/1026
5	I	0.48	1/7778 (0.0%)	0.87	6/11992 (0.1%)
6	J	0.49	2/7770 (0.0%)	0.89	5/11998 (0.0%)
7	S	0.58	0/682	0.80	2/911 (0.2%)
All	All	0.55	3/28869 (0.0%)	0.88	41/41839 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	I	1	DA	OP3-P	-9.95	1.49	1.61
6	J	1	DA	OP3-P	-9.41	1.49	1.61
6	J	113	DA	O3'-P	-5.64	1.54	1.61

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	95	ARG	NE-CZ-NH1	9.26	124.93	120.30
1	E	69	ARG	NE-CZ-NH1	7.09	123.84	120.30
1	E	129	ARG	NE-CZ-NH1	7.08	123.84	120.30
2	B	78	ARG	NE-CZ-NH1	7.08	123.84	120.30
1	O	129	ARG	NE-CZ-NH1	6.90	123.75	120.30
3	Q	29	ARG	NE-CZ-NH1	6.76	123.68	120.30
5	I	207	DT	C1'-O4'-C4'	-6.52	103.58	110.10
1	E	69	ARG	NE-CZ-NH2	-6.50	117.05	120.30
1	E	63	ARG	NE-CZ-NH1	6.49	123.54	120.30
1	A	69	ARG	NE-CZ-NH1	6.31	123.45	120.30
3	M	71	ARG	NE-CZ-NH1	6.23	123.42	120.30
3	M	81	ARG	NE-CZ-NH1	6.22	123.41	120.30
5	I	168	DG	C1'-O4'-C4'	-6.03	104.08	110.10
1	K	131	ARG	NE-CZ-NH1	5.83	123.22	120.30
6	J	101	DA	O5'-P-OP2	-5.83	100.45	105.70
5	I	168	DG	N9-C1'-C2'	5.83	123.67	112.60
6	J	283	DG	O5'-P-OP1	-5.80	100.48	105.70
2	L	78	ARG	NE-CZ-NH1	5.79	123.20	120.30
5	I	331	DA	C1'-O4'-C4'	-5.78	104.32	110.10
5	I	329	DA	C1'-O4'-C4'	-5.67	104.43	110.10
3	C	71	ARG	NE-CZ-NH1	5.66	123.13	120.30
1	O	52	ARG	NE-CZ-NH2	-5.63	117.48	120.30
3	C	81	ARG	NE-CZ-NH1	5.55	123.08	120.30
1	O	69	ARG	NE-CZ-NH1	5.54	123.07	120.30
6	J	106	DG	C1'-O4'-C4'	-5.49	104.61	110.10
6	J	174	DC	C1'-O4'-C4'	-5.46	104.64	110.10
7	S	121	LEU	CA-CB-CG	5.44	127.80	115.30
2	P	10	LEU	CA-CB-CG	5.39	127.70	115.30
3	M	32	ARG	NE-CZ-NH2	5.37	122.98	120.30
4	N	106	LEU	CA-CB-CG	5.32	127.53	115.30
7	S	57	ARG	NE-CZ-NH1	5.26	122.93	120.30
3	Q	42	ARG	NE-CZ-NH1	5.25	122.93	120.30
4	H	80	LEU	CA-CB-CG	5.25	127.37	115.30
5	I	45	DG	C1'-O4'-C4'	-5.23	104.87	110.10
4	H	99	ARG	NE-CZ-NH1	5.12	122.86	120.30
6	J	293	DA	O5'-P-OP2	5.12	116.84	110.70
3	Q	29	ARG	NE-CZ-NH2	-5.08	117.76	120.30
3	C	77	ARG	NE-CZ-NH1	5.07	122.84	120.30
1	E	131	ARG	NE-CZ-NH1	5.04	122.82	120.30
2	F	78	ARG	NE-CZ-NH1	5.04	122.82	120.30
1	K	83	ARG	NE-CZ-NH1	5.02	122.81	120.30

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	1	0
2	B	673	0	722	2	0
2	F	703	0	755	3	0
2	L	683	0	729	9	0
2	P	741	0	796	4	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	9	0
4	N	766	0	797	7	0
4	R	755	0	784	11	0
5	I	6923	0	3777	39	0
6	J	6937	0	3817	50	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
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

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
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	0	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	145	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 (145) 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

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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
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

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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: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
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

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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
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
6:J:178:DT:C6	6:J:179:DT:H72	2.55	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
2:L:101:GLY:O	2:L:102:GLY:C	2.59	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

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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
4:R:70:PHE:CD1	4:R:70:PHE:C	2.94	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
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%)	13	24
2	F	85/103 (82%)	79 (93%)	4 (5%)	2 (2%)	6	9
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%)	15	28

Continued on next page...

*Continued from previous page...*

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	M	109/130 (84%)	103 (94%)	5 (5%)	1 (1%)	17	31
3	Q	104/130 (80%)	102 (98%)	2 (2%)	0	100	100
4	D	94/126 (75%)	93 (99%)	0	1 (1%)	14	26
4	H	94/126 (75%)	90 (96%)	4 (4%)	0	100	100
4	N	95/126 (75%)	90 (95%)	4 (4%)	1 (1%)	14	26
4	R	94/126 (75%)	91 (97%)	2 (2%)	1 (1%)	14	26
7	S	82/213 (38%)	72 (88%)	8 (10%)	2 (2%)	6	9
All	All	1618/2193 (74%)	1558 (96%)	50 (3%)	10 (1%)	25	43

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%)	84 (99%)	1 (1%)	71	88
1	E	87/111 (78%)	81 (93%)	6 (7%)	15	30
1	K	86/111 (78%)	84 (98%)	2 (2%)	50	76
1	O	87/111 (78%)	83 (95%)	4 (5%)	27	50
2	B	69/79 (87%)	69 (100%)	0	100	100

*Continued on next page...*

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	F	72/79 (91%)	70 (97%)	2 (3%)	43	70
2	L	70/79 (89%)	68 (97%)	2 (3%)	42	69
2	P	74/79 (94%)	67 (90%)	7 (10%)	8	17
3	C	88/100 (88%)	85 (97%)	3 (3%)	37	63
3	G	85/100 (85%)	79 (93%)	6 (7%)	14	28
3	M	88/100 (88%)	86 (98%)	2 (2%)	50	76
3	Q	84/100 (84%)	77 (92%)	7 (8%)	11	22
4	D	82/105 (78%)	79 (96%)	3 (4%)	34	60
4	H	82/105 (78%)	73 (89%)	9 (11%)	6	12
4	N	83/105 (79%)	76 (92%)	7 (8%)	11	21
4	R	82/105 (78%)	77 (94%)	5 (6%)	18	36
7	S	73/157 (46%)	58 (80%)	15 (20%)	1	2
All	All	1377/1737 (79%)	1296 (94%)	81 (6%)	19	37

All (81) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	122	LYS
3	C	118	LYS
3	C	119	LYS
3	C	120	THR
4	D	34	LYS
4	D	46	LYS
4	D	86	ARG
1	E	37	LYS
1	E	48	LEU
1	E	53	ARG
1	E	63	ARG
1	E	129	ARG
1	E	134	ARG
2	F	22	LEU
2	F	27	GLN
3	G	29	ARG
3	G	36	LYS
3	G	73	ASN
3	G	74	LYS
3	G	81	ARG
3	G	118	LYS

Continued on next page...

*Continued from previous page...*

Mol	Chain	Res	Type
4	H	31	ARG
4	H	32	SER
4	H	36	SER
4	H	42	TYR
4	H	48	VAL
4	H	52	THR
4	H	80	LEU
4	H	86	ARG
4	H	125	LYS
1	K	63	ARG
1	K	129	ARG
2	L	19	ARG
2	L	95	ARG
3	M	24	GLN
3	M	74	LYS
4	N	31	ARG
4	N	32	SER
4	N	36	SER
4	N	63	ASN
4	N	85	LYS
4	N	86	ARG
4	N	106	LEU
1	O	48	LEU
1	O	117	VAL
1	O	129	ARG
1	O	134	ARG
2	P	10	LEU
2	P	12	LYS
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
4	R	46	LYS
4	R	52	THR

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
4	R	80	LEU
4	R	85	LYS
4	R	125	LYS
7	S	42	LYS
7	S	44	GLN
7	S	51	LEU
7	S	57	ARG
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	113	SER
7	S	116	LEU
7	S	121	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	76	GLN
4	D	95	GLN
1	E	39	HIS
1	E	76	GLN
3	G	73	ASN
1	K	39	HIS
3	M	31	HIS
2	P	27	GLN
2	P	64	ASN
7	S	38	ASN
7	S	44	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 [i](#)

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

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

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 [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	98/136 (72%)	0.34	3 (3%) 49 52	45, 59, 90, 110	0
1	E	100/136 (73%)	0.20	0 100 100	39, 53, 88, 133	0
1	K	99/136 (72%)	0.29	4 (4%) 38 41	33, 46, 86, 130	0
1	O	100/136 (73%)	0.34	2 (2%) 65 68	32, 44, 76, 127	0
2	B	84/103 (81%)	0.59	5 (5%) 21 22	42, 55, 124, 161	0
2	F	87/103 (84%)	0.49	5 (5%) 23 25	39, 50, 137, 162	0
2	L	85/103 (82%)	0.64	6 (7%) 16 16	36, 46, 126, 174	0
2	P	94/103 (91%)	1.22	15 (15%) 1 1	34, 44, 157, 174	0
3	C	112/130 (86%)	0.43	5 (4%) 33 36	41, 53, 115, 151	0
3	G	107/130 (82%)	0.36	6 (5%) 24 25	44, 57, 96, 132	0
3	M	111/130 (85%)	0.42	2 (1%) 68 71	36, 45, 97, 144	0
3	Q	106/130 (81%)	0.20	2 (1%) 66 69	37, 50, 81, 128	0
4	D	96/126 (76%)	0.36	3 (3%) 49 52	43, 55, 109, 129	0
4	H	96/126 (76%)	0.36	6 (6%) 20 21	45, 57, 99, 134	0
4	N	97/126 (76%)	0.33	3 (3%) 49 52	34, 47, 96, 144	0
4	R	96/126 (76%)	0.35	2 (2%) 63 66	38, 49, 93, 127	0
5	I	338/338 (100%)	0.03	25 (7%) 14 15	55, 94, 198, 230	0
6	J	338/338 (100%)	0.02	19 (5%) 24 25	53, 96, 204, 251	0
7	S	84/213 (39%)	4.12	71 (84%) 0 0	112, 147, 171, 188	0
All	All	2328/2869 (81%)	0.44	184 (7%) 12 12	32, 60, 154, 251	0

All (184) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
7	S	39	SER	12.8

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
2	P	15	ALA	10.3
2	B	22	LEU	10.0
7	S	77	VAL	10.0
2	P	17	ARG	9.7
2	B	21	VAL	9.6
2	L	19	ARG	9.6
2	P	10	LEU	9.2
2	P	18	HIS	9.2
7	S	43	ASN	9.1
7	S	121	LEU	8.8
2	B	19	ARG	8.7
3	Q	14	ALA	8.2
7	S	51	LEU	8.1
2	L	18	HIS	8.0
7	S	114	PHE	8.0
7	S	40	LYS	7.7
7	S	102	LEU	7.5
2	L	20	LYS	7.2
2	L	21	VAL	7.2
7	S	38	ASN	7.1
3	C	10	ALA	7.1
7	S	79	TRP	7.1
2	P	16	LYS	6.4
7	S	103	LEU	6.2
7	S	105	VAL	6.2
7	S	86	ARG	6.1
7	S	109	GLY	6.1
4	N	30	LYS	6.1
3	G	13	LYS	6.0
7	S	120	LYS	6.0
7	S	113	SER	6.0
3	G	14	ALA	6.0
7	S	78	PRO	5.9
7	S	117	ASN	5.9
7	S	52	VAL	5.8
7	S	63	ASN	5.8
2	P	19	ARG	5.8
2	B	20	LYS	5.8
1	A	38	PRO	5.8
7	S	82	GLN	5.7
1	K	38	PRO	5.7
7	S	108	THR	5.7

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
5	I	5	DC	5.5
2	F	18	HIS	5.1
3	M	120	THR	5.1
7	S	67	LEU	5.1
7	S	93	ILE	4.9
2	P	21	VAL	4.9
5	I	6	DG	4.9
7	S	116	LEU	4.9
7	S	55	THR	4.8
2	B	23	ARG	4.8
2	L	22	LEU	4.8
7	S	41	LYS	4.8
7	S	119	LYS	4.7
3	G	118	LYS	4.6
6	J	336	DG	4.6
7	S	74	ALA	4.6
7	S	71	TYR	4.5
7	S	100	ASP	4.5
7	S	42	LYS	4.5
2	F	21	VAL	4.5
3	G	119	LYS	4.5
7	S	76	LYS	4.4
6	J	179	DT	4.4
7	S	101	THR	4.4
7	S	59	LEU	4.4
4	H	124	ALA	4.3
7	S	99	ASN	4.3
3	Q	15	LYS	4.2
7	S	95	ALA	4.2
7	S	88	TYR	4.2
4	D	125	LYS	4.1
7	S	96	LEU	4.1
7	S	118	ARG	4.1
2	P	9	GLY	4.1
7	S	97	VAL	4.1
4	H	125	LYS	4.1
2	P	14	GLY	4.0
4	R	125	LYS	3.9
6	J	5	DC	3.9
5	I	7	DG	3.9
4	H	31	ARG	3.9
6	J	185	DC	3.8

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
7	S	44	GLN	3.8
2	P	12	LYS	3.7
5	I	164	DA	3.7
2	P	13	GLY	3.7
5	I	158	DA	3.7
6	J	178	DT	3.7
7	S	54	GLU	3.6
7	S	56	ILE	3.6
5	I	336	DG	3.6
1	O	36	LYS	3.6
7	S	87	THR	3.6
3	C	12	ALA	3.5
2	F	19	ARG	3.5
4	N	29	ARG	3.5
2	L	23	ARG	3.5
3	G	15	LYS	3.5
5	I	335	DC	3.4
5	I	4	DG	3.3
6	J	4	DG	3.3
2	P	22	LEU	3.3
7	S	80	PHE	3.3
7	S	107	GLY	3.3
7	S	65	SER	3.2
7	S	48	TYR	3.2
7	S	64	GLY	3.2
2	P	11	GLY	3.2
7	S	115	LYS	3.2
6	J	166	DC	3.1
3	C	11	ARG	3.1
5	I	328	DA	3.1
4	D	31	ARG	3.1
7	S	112	GLY	3.1
2	F	20	LYS	3.1
7	S	70	ILE	3.1
7	S	90	LYS	3.1
6	J	337	DA	3.1
6	J	6	DG	3.1
5	I	327	DA	3.0
4	D	32	SER	3.0
7	S	68	ALA	2.9
2	F	17	ARG	2.9
6	J	3	DC	2.9

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
7	S	75	LYS	2.9
5	I	338	DT	2.9
7	S	91	TYR	2.9
7	S	104	GLN	2.8
3	C	13	LYS	2.8
7	S	92	SER	2.8
3	M	121	GLU	2.8
5	I	337	DA	2.8
6	J	335	DC	2.8
1	A	135	ALA	2.8
7	S	84	ASN	2.8
5	I	161	DA	2.8
6	J	164	DT	2.7
3	C	120	THR	2.7
6	J	175	DT	2.7
7	S	45	PRO	2.6
3	G	36	LYS	2.6
7	S	50	GLN	2.6
2	P	20	LYS	2.6
5	I	162	DA	2.6
6	J	173	DG	2.6
7	S	62	ARG	2.6
7	S	53	VAL	2.5
1	K	39	HIS	2.5
5	I	3	DC	2.5
7	S	72	THR	2.5
7	S	98	GLN	2.5
7	S	57	ARG	2.5
2	P	23	ARG	2.5
1	O	37	LYS	2.5
5	I	157	DA	2.4
5	I	168	DG	2.4
5	I	175	DA	2.4
7	S	58	ARG	2.4
5	I	165	DG	2.4
7	S	81	ASP	2.3
5	I	160	DA	2.3
5	I	177	DA	2.3
6	J	174	DC	2.3
5	I	159	DA	2.3
4	H	34	LYS	2.3
1	A	39	HIS	2.2

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
5	I	106	DC	2.2
5	I	172	DC	2.2
1	K	37	LYS	2.2
4	H	33	ARG	2.2
4	R	82	HIS	2.2
6	J	167	DG	2.2
7	S	66	SER	2.1
5	I	173	DG	2.1
6	J	165	DG	2.1
4	N	31	ARG	2.1
6	J	212	DG	2.0
4	H	42	TYR	2.0
1	K	49	ARG	2.0
6	J	1	DA	2.0
7	S	111	ASN	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 monosaccharides 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
9	CA	J	405	1/1	0.69	0.18	112,112,112,112	0
9	CA	G	201	1/1	0.72	0.12	97,97,97,97	0
9	CA	I	420	1/1	0.74	0.10	108,108,108,108	0
9	CA	J	410	1/1	0.75	0.10	103,103,103,103	0
10	K	I	433	1/1	0.76	0.11	114,114,114,114	0
9	CA	I	427	1/1	0.78	0.15	102,102,102,102	0
9	CA	I	419	1/1	0.79	0.13	109,109,109,109	0
9	CA	J	422	1/1	0.80	0.17	95,95,95,95	0

*Continued on next page...*

*Continued from previous page...*

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
9	CA	I	426	1/1	0.81	0.11	104,104,104,104	0
9	CA	I	402	1/1	0.81	0.10	103,103,103,103	0
9	CA	I	405	1/1	0.81	0.12	102,102,102,102	0
9	CA	I	401	1/1	0.82	0.10	110,110,110,110	0
9	CA	I	417	1/1	0.82	0.11	108,108,108,108	0
9	CA	J	408	1/1	0.83	0.09	104,104,104,104	0
9	CA	R	201	1/1	0.83	0.15	97,97,97,97	0
9	CA	J	411	1/1	0.83	0.08	94,94,94,94	0
9	CA	I	416	1/1	0.84	0.16	115,115,115,115	0
9	CA	I	415	1/1	0.84	0.12	102,102,102,102	0
10	K	C	205	1/1	0.84	0.13	93,93,93,93	0
9	CA	J	413	1/1	0.84	0.10	106,106,106,106	0
10	K	I	434	1/1	0.85	0.09	110,110,110,110	0
9	CA	C	203	1/1	0.86	0.10	87,87,87,87	0
9	CA	I	425	1/1	0.86	0.21	104,104,104,104	0
9	CA	J	412	1/1	0.87	0.08	90,90,90,90	0
9	CA	M	202	1/1	0.87	0.19	94,94,94,94	0
9	CA	I	407	1/1	0.87	0.05	96,96,96,96	0
10	K	I	432	1/1	0.88	0.05	99,99,99,99	0
9	CA	J	421	1/1	0.88	0.10	100,100,100,100	0
10	K	I	429	1/1	0.88	0.10	80,80,80,80	0
9	CA	I	418	1/1	0.90	0.09	97,97,97,97	0
9	CA	G	203	1/1	0.90	0.10	102,102,102,102	0
9	CA	I	406	1/1	0.90	0.04	106,106,106,106	0
9	CA	I	428	1/1	0.90	0.32	82,82,82,82	0
9	CA	J	402	1/1	0.90	0.29	111,111,111,111	0
9	CA	I	424	1/1	0.90	0.07	100,100,100,100	0
9	CA	J	406	1/1	0.90	0.08	83,83,83,83	0
9	CA	J	423	1/1	0.91	0.09	100,100,100,100	0
10	K	I	431	1/1	0.91	0.06	97,97,97,97	0
9	CA	I	403	1/1	0.91	0.06	102,102,102,102	0
9	CA	I	409	1/1	0.91	0.14	85,85,85,85	0
9	CA	I	421	1/1	0.91	0.04	102,102,102,102	0
9	CA	I	422	1/1	0.92	0.08	91,91,91,91	0
9	CA	N	201	1/1	0.92	0.14	84,84,84,84	0
9	CA	Q	202	1/1	0.92	0.08	91,91,91,91	0
9	CA	I	404	1/1	0.92	0.06	86,86,86,86	0
9	CA	J	416	1/1	0.92	0.07	95,95,95,95	0
9	CA	J	404	1/1	0.93	0.05	93,93,93,93	0
9	CA	I	413	1/1	0.93	0.12	98,98,98,98	0
9	CA	I	414	1/1	0.93	0.27	111,111,111,111	0
9	CA	C	202	1/1	0.93	0.07	96,96,96,96	0

*Continued on next page...*



*Continued from previous page...*

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
9	CA	J	419	1/1	0.93	0.08	101,101,101,101	0
8	CL	A	201	1/1	0.93	0.15	72,72,72,72	0
9	CA	J	409	1/1	0.94	0.08	102,102,102,102	0
10	K	I	430	1/1	0.94	0.18	80,80,80,80	0
9	CA	J	420	1/1	0.94	0.07	111,111,111,111	0
9	CA	J	415	1/1	0.94	0.10	82,82,82,82	0
9	CA	I	423	1/1	0.94	0.10	95,95,95,95	0
9	CA	J	417	1/1	0.94	0.12	85,85,85,85	0
9	CA	I	411	1/1	0.95	0.05	83,83,83,83	0
9	CA	I	412	1/1	0.95	0.06	92,92,92,92	0
9	CA	J	418	1/1	0.95	0.16	101,101,101,101	0
9	CA	J	407	1/1	0.95	0.04	99,99,99,99	0
9	CA	Q	201	1/1	0.95	0.09	98,98,98,98	0
9	CA	I	408	1/1	0.95	0.04	91,91,91,91	0
8	CL	K	201	1/1	0.95	0.12	64,64,64,64	0
10	K	J	424	1/1	0.95	0.07	77,77,77,77	0
10	K	M	204	1/1	0.95	0.07	80,80,80,80	0
9	CA	C	201	1/1	0.96	0.17	87,87,87,87	0
9	CA	G	202	1/1	0.96	0.08	117,117,117,117	0
9	CA	N	202	1/1	0.96	0.20	83,83,83,83	0
9	CA	J	414	1/1	0.96	0.10	95,95,95,95	0
10	K	J	425	1/1	0.96	0.12	77,77,77,77	0
9	CA	I	410	1/1	0.96	0.04	85,85,85,85	0
8	CL	E	201	1/1	0.97	0.07	66,66,66,66	0
9	CA	D	201	1/1	0.97	0.15	80,80,80,80	0
8	CL	O	201	1/1	0.97	0.10	49,49,49,49	0
9	CA	M	201	1/1	0.97	0.24	91,91,91,91	0
9	CA	J	401	1/1	0.98	0.16	88,88,88,88	0
8	CL	Q	203	1/1	0.98	0.09	51,51,51,51	0
9	CA	J	403	1/1	0.98	0.13	79,79,79,79	0
8	CL	G	204	1/1	0.98	0.14	61,61,61,61	0
8	CL	C	204	1/1	0.99	0.09	55,55,55,55	0
8	CL	M	203	1/1	1.00	0.12	48,48,48,48	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.