



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:57 PM UTC

PDB ID : 6L9M / pdb_00006l9m
Title : H2-Ld complexed with AH1 peptide
Authors : Wei, P.C.; Yin, L.
Deposited on : 2019-11-10
Resolution : 2.60 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

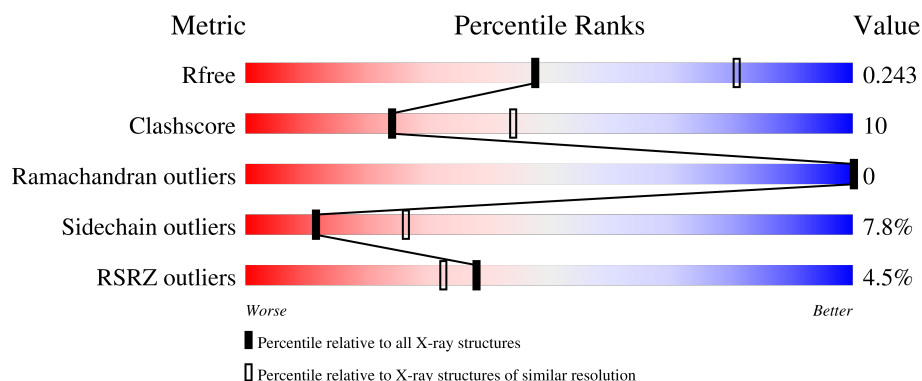
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	278	5% 80% 18% .
1	D	278	8% 78% 18% .
1	G	278	4% 77% 21% .
1	J	278	4% 79% 18% .
2	B	99	4% 72% 24% .

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Mol	Chain	Length	Quality of chain
2	E	99	2%68%27%••
2	H	99	4%69%28%•
2	K	99	3%71%24%5%
3	C	9	89%11%
3	F	9	22%89%11%
3	I	9	11%78%22%
3	L	9	78%22%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 12891 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 H2-Ld.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	278	Total	C	N	O	S	41	0	0
			2268	1437	395	426	10			
1	D	278	Total	C	N	O	S	41	0	0
			2268	1437	395	426	10			
1	G	278	Total	C	N	O	S	41	0	0
			2268	1437	395	426	10			
1	J	278	Total	C	N	O	S	41	0	0
			2268	1437	395	426	10			

- Molecule 2 is a protein called b2m.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	99	Total	C	N	O	S	10	0	0
			821	524	138	152	7			
2	E	99	Total	C	N	O	S	10	0	0
			821	524	138	152	7			
2	H	99	Total	C	N	O	S	10	0	0
			821	524	138	152	7			
2	K	99	Total	C	N	O	S	10	0	0
			821	524	138	152	7			

- Molecule 3 is a protein called SER-PRO-SER-TYR-VAL-TYR-HIS-GLN-PHE.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	9	Total	C	N	O	0	0	0
			81	54	12	15			
3	F	9	Total	C	N	O	0	0	0
			81	54	12	15			
3	I	9	Total	C	N	O	0	0	0
			81	54	12	15			
3	L	9	Total	C	N	O	0	0	0
			81	54	12	15			

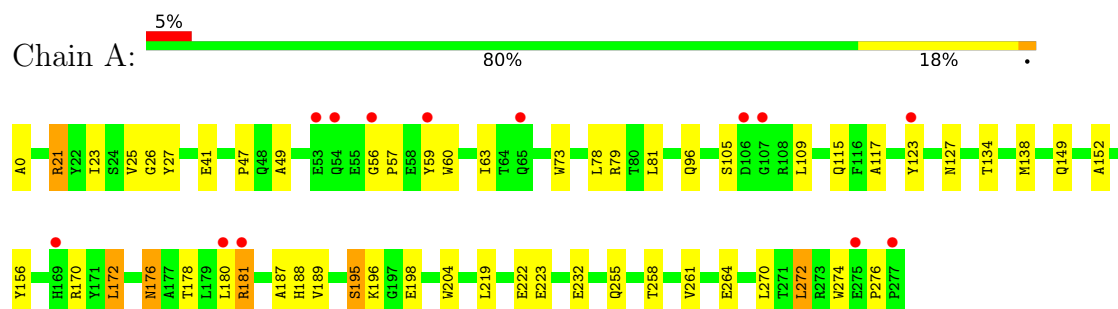
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	32	Total 32	O 32	0	0
4	B	12	Total 12	O 12	0	0
4	D	36	Total 36	O 36	0	0
4	E	21	Total 21	O 21	0	0
4	G	38	Total 38	O 38	0	0
4	H	19	Total 19	O 19	0	0
4	J	32	Total 32	O 32	0	0
4	K	19	Total 19	O 19	0	0
4	L	2	Total 2	O 2	0	0

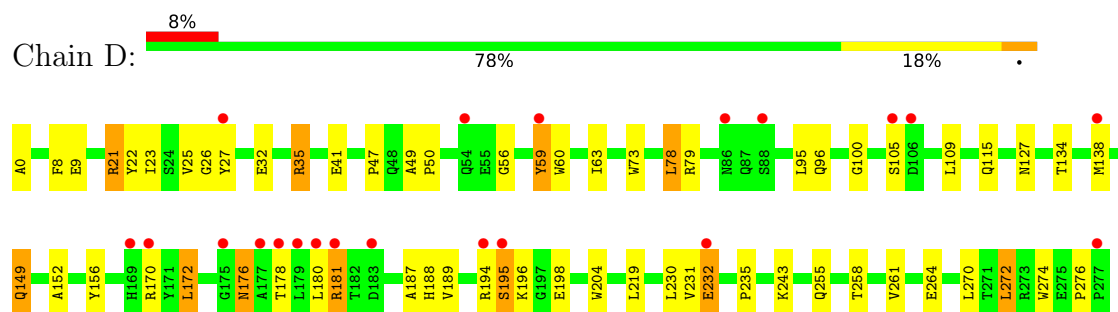
3 Residue-property plots [i](#)

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.

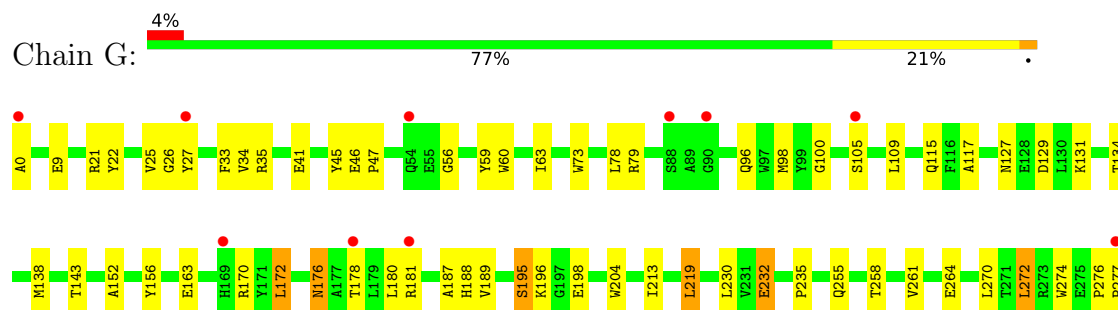
• Molecule 1: H2-Ld



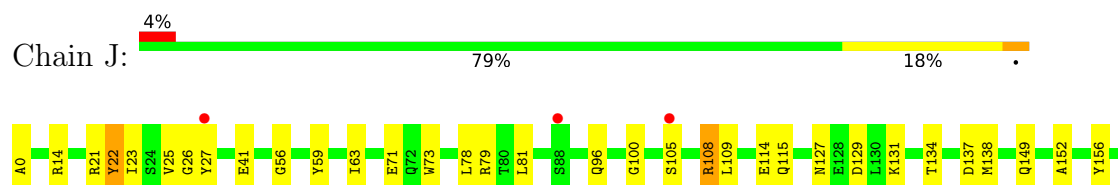
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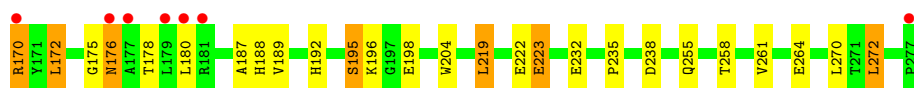


• Molecule 1: H2-Ld

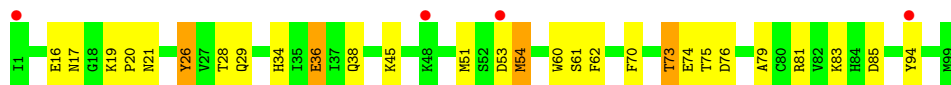
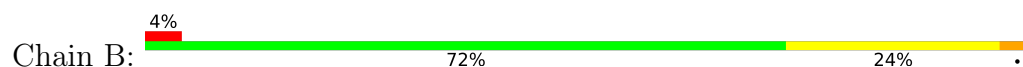


• Molecule 1: H2-Ld





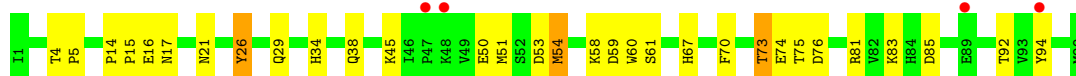
- Molecule 2: b2m



- Molecule 2: b2m



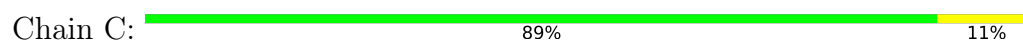
- Molecule 2: b2m



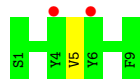
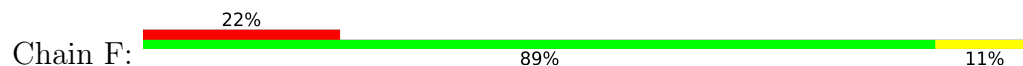
- Molecule 2: b2m



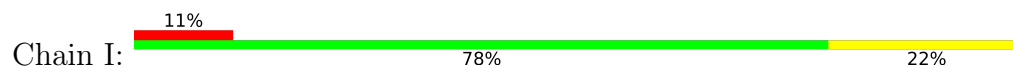
- Molecule 3: SER-PRO-SER-TYR-VAL-TYR-HIS-GLN-PHE



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Chain L: 78% 22%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	51.87Å 88.39Å 105.71Å 80.97° 75.96° 88.24°	Depositor
Resolution (Å)	19.97 – 2.60 19.97 – 2.60	Depositor EDS
% Data completeness (in resolution range)	95.7 (19.97-2.60) 95.5 (19.97-2.60)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.08 (at 2.59Å)	Xtriage
Refinement program	PHENIX 1.7.3_928	Depositor
R, R_{free}	0.213 , 0.249 0.204 , 0.243	Depositor DCC
R_{free} test set	2006 reflections (3.38%)	wwPDB-VP
Wilson B-factor (Å ²)	29.5	Xtriage
Anisotropy	0.539	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 38.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.018 for h,-k,h-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	12891	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.09% 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

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	0.62	2/2337 (0.1%)	0.91	5/3180 (0.2%)
1	D	0.68	5/2337 (0.2%)	0.97	10/3180 (0.3%)
1	G	0.65	2/2337 (0.1%)	0.90	5/3180 (0.2%)
1	J	0.83	9/2337 (0.4%)	0.99	12/3180 (0.4%)
2	B	0.67	4/847 (0.5%)	0.92	2/1148 (0.2%)
2	E	0.97	7/847 (0.8%)	1.01	5/1148 (0.4%)
2	H	0.69	4/847 (0.5%)	0.88	1/1148 (0.1%)
2	K	0.88	9/847 (1.1%)	0.92	1/1148 (0.1%)
3	C	0.40	0/85	0.64	0/114
3	F	0.37	0/85	0.73	0/114
3	I	0.41	0/85	0.69	0/114
3	L	0.44	0/85	0.83	0/114
All	All	0.72	42/13076 (0.3%)	0.94	41/17768 (0.2%)

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	E	0	1

All (42) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	33	PRO	C-O	-14.95	1.03	1.24
1	D	35	ARG	CZ-NH2	-10.85	1.19	1.33
1	J	22	TYR	CG-CD2	-10.62	1.17	1.39
1	J	22	TYR	CG-CD1	-10.37	1.17	1.39
1	J	22	TYR	CE2-CZ	-9.14	1.16	1.38
1	J	22	TYR	CE1-CZ	-8.98	1.16	1.38
2	E	34	HIS	CD2-NE2	-8.56	1.28	1.37
1	J	223	GLU	CD-OE2	-8.53	1.09	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	J	223	GLU	CD-OE1	-8.02	1.10	1.25
1	J	175	GLY	C-O	-7.36	1.19	1.24
2	K	78	TYR	CG-CD1	-7.13	1.24	1.39
2	E	26	TYR	CE1-CZ	-6.56	1.22	1.38
2	K	26	TYR	CE1-CZ	-6.48	1.22	1.38
2	E	26	TYR	CG-CD1	-6.44	1.25	1.39
2	E	26	TYR	CE2-CZ	-6.38	1.23	1.38
2	K	26	TYR	CG-CD1	-6.36	1.25	1.39
2	E	26	TYR	CG-CD2	-6.35	1.26	1.39
2	K	26	TYR	CG-CD2	-6.33	1.26	1.39
2	B	26	TYR	CE1-CZ	-6.30	1.23	1.38
2	K	78	TYR	CE1-CZ	-6.27	1.23	1.38
2	H	26	TYR	CG-CD1	-6.25	1.26	1.39
2	H	26	TYR	CE1-CZ	-6.25	1.23	1.38
2	K	78	TYR	CG-CD2	-6.22	1.26	1.39
2	B	26	TYR	CG-CD2	-6.17	1.26	1.39
2	B	26	TYR	CG-CD1	-6.15	1.26	1.39
2	B	26	TYR	CE2-CZ	-6.11	1.23	1.38
2	K	26	TYR	CE2-CZ	-6.00	1.23	1.38
2	H	26	TYR	CG-CD2	-5.97	1.26	1.39
2	K	78	TYR	CE2-CZ	-5.97	1.24	1.38
1	D	181	ARG	CD-NE	-5.96	1.38	1.46
2	H	26	TYR	CE2-CZ	-5.95	1.24	1.38
1	A	219	LEU	CG-CD1	-5.91	1.33	1.52
1	D	149	GLN	CD-OE1	-5.89	1.12	1.23
1	D	219	LEU	CG-CD2	-5.86	1.33	1.52
1	A	219	LEU	CG-CD2	-5.76	1.33	1.52
1	D	219	LEU	CG-CD1	-5.66	1.33	1.52
2	K	89	GLU	CD-OE1	-5.66	1.14	1.25
1	J	219	LEU	CG-CD2	-5.60	1.34	1.52
1	G	219	LEU	CG-CD2	-5.49	1.34	1.52
1	J	219	LEU	CG-CD1	-5.30	1.35	1.52
1	G	181	ARG	CG-CD	5.30	1.68	1.52
2	E	34	HIS	CG-ND1	5.26	1.44	1.38

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	35	ARG	NE-CZ-NH1	15.52	137.02	121.50
1	J	219	LEU	CD1-CG-CD2	-11.04	86.52	110.80
1	A	219	LEU	CD1-CG-CD2	-10.85	86.94	110.80
1	G	219	LEU	CD1-CG-CD2	-10.84	86.96	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	219	LEU	CD1-CG-CD2	-10.74	87.17	110.80
2	E	34	HIS	CB-CA-C	8.92	123.81	109.90
2	E	34	HIS	CA-C-O	-8.76	110.94	120.92
2	B	36	GLU	CA-CB-CG	8.56	131.22	114.10
1	J	223	GLU	OE1-CD-OE2	-8.50	102.50	122.90
1	J	170	ARG	CB-CG-CD	7.82	129.28	111.30
1	J	79	ARG	NE-CZ-NH2	7.63	126.07	119.20
1	D	35	ARG	NH1-CZ-NH2	-7.15	110.01	119.30
1	D	35	ARG	NE-CZ-NH2	-7.04	112.86	119.20
1	J	170	ARG	CB-CA-C	-6.97	99.83	110.92
1	J	79	ARG	NH1-CZ-NH2	-6.60	110.72	119.30
2	E	34	HIS	ND1-CE1-NE2	6.57	114.97	108.40
1	D	149	GLN	CB-CG-CD	6.43	123.53	112.60
1	A	56	GLY	CA-C-N	6.31	126.22	119.28
1	A	56	GLY	C-N-CA	6.31	126.22	119.28
1	G	56	GLY	CA-C-N	6.18	126.08	119.28
1	G	56	GLY	C-N-CA	6.18	126.08	119.28
1	J	108	ARG	NE-CZ-NH2	6.08	124.67	119.20
1	D	181	ARG	CG-CD-NE	-6.05	98.69	112.00
1	J	22	TYR	CE1-CZ-CE2	-6.04	108.21	120.30
1	D	56	GLY	CA-C-N	5.99	125.86	119.28
1	D	56	GLY	C-N-CA	5.99	125.86	119.28
1	J	56	GLY	CA-C-N	5.91	125.79	119.28
1	J	56	GLY	C-N-CA	5.91	125.79	119.28
2	E	34	HIS	CG-ND1-CE1	-5.88	99.30	109.30
2	E	59	ASP	N-CA-C	-5.54	102.28	110.48
1	G	181	ARG	CG-CD-NE	-5.51	99.88	112.00
2	H	59	ASP	N-CA-C	-5.47	101.56	110.20
2	K	59	ASP	N-CA-C	-5.42	102.45	110.48
1	D	100	GLY	N-CA-C	5.24	117.09	110.38
1	J	100	GLY	N-CA-C	5.19	117.02	110.38
1	A	49	ALA	CA-C-N	5.12	124.91	119.28
1	A	49	ALA	C-N-CA	5.12	124.91	119.28
2	B	36	GLU	N-CA-C	5.11	117.30	109.07
1	D	35	ARG	CD-NE-CZ	-5.10	117.26	124.40
1	J	170	ARG	CA-CB-CG	5.08	124.26	114.10
1	G	100	GLY	N-CA-C	5.06	116.86	110.38

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	E	34	HIS	Mainchain

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	2268	0	2131	45	0
1	D	2268	0	2131	56	0
1	G	2268	0	2131	42	0
1	J	2268	0	2131	40	0
2	B	821	0	796	18	0
2	E	821	0	796	21	0
2	H	821	0	796	27	0
2	K	821	0	796	28	0
3	C	81	0	70	1	0
3	F	81	0	70	1	0
3	I	81	0	70	3	0
3	L	81	0	70	1	0
4	A	32	0	0	0	0
4	B	12	0	0	0	0
4	D	36	0	0	2	0
4	E	21	0	0	0	0
4	G	38	0	0	4	0
4	H	19	0	0	0	0
4	J	32	0	0	3	0
4	K	19	0	0	1	0
4	L	2	0	0	0	0
All	All	12891	0	11988	241	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 10.

All (241) 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)
2:K:70:PHE:HD2	2:K:78:TYR:CE1	1.74	1.04
1:A:96:GLN:HE22	2:B:60:TRP:HB3	1.18	1.02

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:59:TYR:HE1	1:D:63:ILE:HD13	1.22	1.01
2:K:70:PHE:CD2	2:K:78:TYR:HE1	1.84	0.94
1:D:59:TYR:CE1	1:D:63:ILE:HD13	2.02	0.94
2:K:70:PHE:HD2	2:K:78:TYR:HE1	0.91	0.90
1:A:57:PRO:HA	1:D:181:ARG:CZ	2.04	0.87
1:A:60:TRP:CD1	1:D:181:ARG:HH11	1.92	0.86
1:A:21:ARG:HH11	1:A:23:ILE:HD11	1.44	0.82
2:K:70:PHE:CD2	2:K:78:TYR:CE1	2.63	0.80
1:G:96:GLN:HE22	2:H:60:TRP:HB3	1.47	0.79
1:A:60:TRP:HD1	1:D:181:ARG:HH11	1.28	0.78
1:A:96:GLN:NE2	2:B:60:TRP:HB3	1.98	0.78
1:D:21:ARG:HH11	1:D:23:ILE:HD11	1.48	0.77
1:A:60:TRP:HD1	1:D:181:ARG:NH1	1.83	0.76
2:H:50:GLU:HB3	2:H:67:HIS:CD2	2.21	0.76
1:D:9:GLU:OE1	1:D:22:TYR:OH	2.05	0.74
1:G:9:GLU:OE1	1:G:22:TYR:OH	2.06	0.73
1:J:25:VAL:HG12	1:J:27:TYR:HE1	1.53	0.72
2:K:34:HIS:O	4:K:101:HOH:O	2.08	0.70
1:G:25:VAL:HG12	1:G:27:TYR:HE1	1.54	0.69
1:J:22:TYR:HD2	1:J:71:GLU:HG3	1.58	0.69
1:J:114:GLU:CD	1:J:156:TYR:HD2	2.01	0.68
1:A:60:TRP:CD1	1:D:181:ARG:NH1	2.58	0.68
1:G:189:VAL:HG23	1:G:272:LEU:HD12	1.75	0.68
1:A:57:PRO:CG	1:D:181:ARG:NH2	2.57	0.67
1:J:189:VAL:HG23	1:J:272:LEU:HD12	1.76	0.67
1:A:57:PRO:HG3	1:D:181:ARG:NH2	2.10	0.67
2:B:38:GLN:OE1	2:B:81:ARG:NH1	2.25	0.66
1:G:96:GLN:NE2	2:H:60:TRP:HB3	2.11	0.65
1:J:22:TYR:HD2	1:J:71:GLU:CG	2.09	0.65
1:D:96:GLN:HE22	2:E:60:TRP:HB3	1.61	0.65
1:J:152:ALA:O	1:J:156:TYR:HD1	1.79	0.65
1:A:189:VAL:HG23	1:A:272:LEU:HD12	1.78	0.65
1:A:25:VAL:HG12	1:A:27:TYR:CE1	2.32	0.65
2:E:38:GLN:OE1	2:E:81:ARG:NH1	2.26	0.65
1:A:59:TYR:CE1	1:A:63:ILE:HD13	2.32	0.64
1:D:189:VAL:HG23	1:D:272:LEU:HD12	1.79	0.64
1:G:25:VAL:CG1	1:G:27:TYR:HE1	2.10	0.64
1:G:59:TYR:CE1	1:G:63:ILE:HD13	2.32	0.64
2:H:73:THR:HG22	2:H:76:ASP:H	1.61	0.64
1:J:26:GLY:C	1:J:27:TYR:HD1	2.05	0.64
2:K:73:THR:HG22	2:K:76:ASP:H	1.62	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:59:TYR:CE1	1:J:63:ILE:HD13	2.32	0.64
1:D:59:TYR:C	1:D:59:TYR:HD1	2.05	0.64
1:G:26:GLY:C	1:G:27:TYR:HD1	2.04	0.64
1:J:25:VAL:CG1	1:J:27:TYR:HE1	2.10	0.64
2:E:73:THR:HG22	2:E:76:ASP:H	1.63	0.63
1:G:235:PRO:HD3	2:H:26:TYR:CE1	2.33	0.63
2:K:73:THR:HG22	2:K:75:THR:H	1.64	0.62
1:D:25:VAL:HG12	1:D:27:TYR:CE1	2.35	0.62
2:B:73:THR:HG22	2:B:76:ASP:H	1.63	0.62
1:A:21:ARG:NH1	1:A:23:ILE:HD11	2.14	0.62
1:J:96:GLN:HE22	2:K:60:TRP:HB3	1.64	0.62
1:D:21:ARG:NH1	1:D:23:ILE:HD11	2.14	0.61
2:B:79:ALA:HB2	2:B:94:TYR:CD1	2.35	0.61
1:D:59:TYR:C	1:D:59:TYR:CD1	2.78	0.61
1:A:57:PRO:CA	1:D:181:ARG:CZ	2.78	0.61
2:K:38:GLN:OE1	2:K:81:ARG:NH1	2.27	0.60
1:A:25:VAL:HG12	1:A:27:TYR:HE1	1.66	0.59
1:D:195:SER:OG	4:D:301:HOH:O	2.15	0.59
1:A:57:PRO:HG3	1:D:181:ARG:HH21	1.65	0.59
1:G:117:ALA:HB2	2:H:60:TRP:CE2	2.38	0.59
2:B:73:THR:HG22	2:B:75:THR:H	1.67	0.58
2:H:38:GLN:OE1	2:H:81:ARG:NH1	2.28	0.58
1:J:0:ALA:HB2	1:J:264:GLU:OE2	2.04	0.58
1:A:172:LEU:O	1:A:176:ASN:HB3	2.04	0.58
2:K:79:ALA:HB2	2:K:94:TYR:CD1	2.38	0.58
1:D:232:GLU:HG3	1:J:222:GLU:OE1	2.04	0.58
1:A:0:ALA:HB2	1:A:264:GLU:OE2	2.04	0.57
1:G:73:TRP:CZ2	3:I:5:VAL:HG11	2.40	0.57
2:H:73:THR:HG22	2:H:75:THR:H	1.67	0.57
2:B:79:ALA:HB2	2:B:94:TYR:HD1	1.68	0.57
2:E:73:THR:HG22	2:E:75:THR:H	1.69	0.57
1:G:0:ALA:HB2	1:G:264:GLU:OE2	2.05	0.57
1:J:238:ASP:O	4:J:301:HOH:O	2.17	0.57
1:A:57:PRO:CA	1:D:181:ARG:NH2	2.68	0.56
1:A:25:VAL:CG1	1:A:27:TYR:HE1	2.18	0.56
1:G:172:LEU:O	1:G:176:ASN:HB3	2.06	0.56
1:D:25:VAL:HG12	1:D:27:TYR:HE1	1.69	0.56
1:A:222:GLU:OE1	1:G:232:GLU:HG3	2.06	0.56
1:D:0:ALA:HB2	1:D:264:GLU:OE2	2.05	0.55
1:D:194:ARG:NE	4:D:305:HOH:O	2.38	0.55
1:D:172:LEU:O	1:D:176:ASN:HB3	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:230:LEU:HD11	1:J:223:GLU:CD	2.32	0.55
1:G:26:GLY:O	1:G:27:TYR:HD1	1.90	0.54
1:D:26:GLY:O	1:D:27:TYR:HD1	1.90	0.54
1:D:25:VAL:CG1	1:D:27:TYR:HE1	2.20	0.54
1:G:27:TYR:HD2	4:G:334:HOH:O	1.91	0.54
1:J:25:VAL:HG12	1:J:27:TYR:CE1	2.40	0.54
2:K:79:ALA:HB2	2:K:94:TYR:HD1	1.73	0.54
1:A:73:TRP:CZ2	3:C:5:VAL:HG11	2.44	0.53
2:H:92:THR:CG2	2:H:94:TYR:HE1	2.21	0.53
1:A:26:GLY:O	1:A:27:TYR:HD1	1.91	0.53
1:J:26:GLY:O	1:J:27:TYR:HD1	1.92	0.53
1:J:137:ASP:OD2	4:J:302:HOH:O	2.19	0.53
1:D:96:GLN:NE2	2:E:60:TRP:HB3	2.24	0.52
1:G:25:VAL:HG12	1:G:27:TYR:CE1	2.41	0.52
1:J:195:SER:OG	1:J:196:LYS:N	2.41	0.52
1:D:8:PHE:HB3	2:E:56:PHE:CE2	2.45	0.52
1:A:195:SER:OG	1:A:196:LYS:N	2.42	0.52
1:G:188:HIS:CE1	1:G:204:TRP:CD1	2.98	0.51
2:K:77:THR:O	2:K:78:TYR:HD2	1.92	0.51
1:D:195:SER:OG	1:D:196:LYS:N	2.44	0.51
1:G:127:ASN:OD1	1:G:134:THR:OG1	2.25	0.50
1:G:152:ALA:O	1:G:156:TYR:HD1	1.94	0.50
1:G:187:ALA:HA	1:G:204:TRP:O	2.11	0.50
1:A:57:PRO:CB	1:D:181:ARG:NH2	2.74	0.50
1:D:187:ALA:HA	1:D:204:TRP:O	2.12	0.50
1:G:163:GLU:HG2	4:G:310:HOH:O	2.11	0.50
1:A:96:GLN:OE1	2:B:62:PHE:CZ	2.64	0.50
1:A:187:ALA:HA	1:A:204:TRP:O	2.11	0.50
1:J:96:GLN:NE2	2:K:60:TRP:HB3	2.27	0.50
1:J:187:ALA:HA	1:J:204:TRP:O	2.12	0.50
1:D:127:ASN:OD1	1:D:134:THR:OG1	2.26	0.49
1:D:188:HIS:CE1	1:D:204:TRP:CD1	3.00	0.49
1:D:25:VAL:CG1	1:D:27:TYR:CE1	2.96	0.49
1:A:223:GLU:CD	1:G:230:LEU:HD11	2.38	0.49
1:D:152:ALA:O	1:D:156:TYR:HD1	1.96	0.49
1:J:192:HIS:CE1	2:K:98:ASP:HB3	2.48	0.49
1:A:127:ASN:OD1	1:A:134:THR:OG1	2.26	0.49
1:J:172:LEU:O	1:J:176:ASN:HB3	2.12	0.49
1:A:25:VAL:CG1	1:A:27:TYR:CE1	2.94	0.49
1:G:34:VAL:CG1	1:G:45:TYR:CD1	2.96	0.48
1:G:195:SER:OG	1:G:196:LYS:N	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:22:TYR:CD1	1:J:22:TYR:C	2.89	0.48
1:J:235:PRO:HD3	2:K:26:TYR:CE1	2.49	0.48
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.49	0.48
2:H:17:ASN:HB3	2:H:74:GLU:OE2	2.14	0.47
2:H:50:GLU:HB3	2:H:67:HIS:NE2	2.27	0.47
1:J:188:HIS:CE1	1:J:204:TRP:CD1	3.02	0.47
2:B:17:ASN:HB3	2:B:74:GLU:OE2	2.14	0.47
1:G:172:LEU:HD12	1:G:172:LEU:HA	1.78	0.47
1:J:114:GLU:CD	1:J:156:TYR:CD2	2.89	0.47
2:K:70:PHE:CD2	2:K:78:TYR:CD1	3.02	0.47
1:J:127:ASN:OD1	1:J:134:THR:OG1	2.26	0.47
1:A:57:PRO:HA	1:D:181:ARG:NH1	2.30	0.47
2:H:50:GLU:HB3	2:H:67:HIS:HD2	1.78	0.46
1:D:231:VAL:O	1:D:243:LYS:NZ	2.48	0.46
2:K:73:THR:CG2	2:K:75:THR:H	2.29	0.46
2:B:26:TYR:CE2	2:B:28:THR:CG2	2.98	0.46
1:G:277:PRO:HD2	4:G:335:HOH:O	2.15	0.46
1:D:59:TYR:HD1	1:D:59:TYR:O	1.99	0.46
1:G:35:ARG:CZ	2:H:53:ASP:HB2	2.46	0.46
1:J:22:TYR:HD1	1:J:23:ILE:N	2.14	0.45
1:J:26:GLY:C	1:J:27:TYR:CD1	2.92	0.45
1:A:152:ALA:O	1:A:156:TYR:HD1	1.99	0.45
2:E:92:THR:CG2	2:E:94:TYR:HE1	2.30	0.45
2:H:21:ASN:HB3	2:H:70:PHE:CE1	2.51	0.45
1:A:176:ASN:O	1:A:180:LEU:HB2	2.17	0.45
1:D:261:VAL:HB	1:D:270:LEU:HB2	1.98	0.45
1:J:261:VAL:HB	1:J:270:LEU:HB2	1.99	0.45
2:K:17:ASN:HB3	2:K:74:GLU:OE2	2.17	0.45
1:D:49:ALA:HA	1:D:50:PRO:HD3	1.88	0.45
1:G:176:ASN:O	1:G:180:LEU:HB2	2.17	0.45
1:D:73:TRP:CZ2	3:F:5:VAL:HG11	2.52	0.45
1:D:255:GLN:CD	1:D:255:GLN:H	2.24	0.45
2:E:73:THR:CG2	2:E:75:THR:H	2.30	0.45
1:G:261:VAL:HB	1:G:270:LEU:HB2	1.99	0.45
1:G:129:ASP:HB2	1:G:131:LYS:HE3	1.99	0.45
2:B:73:THR:CG2	2:B:75:THR:H	2.30	0.45
2:H:73:THR:CG2	2:H:75:THR:H	2.30	0.44
2:E:54:MET:HE2	2:E:54:MET:HB2	1.85	0.44
1:J:73:TRP:CZ2	3:L:5:VAL:HG11	2.51	0.44
1:J:255:GLN:H	1:J:255:GLN:CD	2.25	0.44
2:K:21:ASN:HB3	2:K:70:PHE:CE1	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:26:GLY:C	1:D:27:TYR:CD1	2.95	0.44
1:D:26:GLY:C	1:D:27:TYR:HD1	2.26	0.44
1:D:176:ASN:O	1:D:180:LEU:HB2	2.18	0.44
1:D:235:PRO:CD	2:E:26:TYR:CE1	3.01	0.44
1:G:26:GLY:C	1:G:27:TYR:CD1	2.90	0.44
1:G:156:TYR:OH	3:I:5:VAL:O	2.35	0.44
1:A:261:VAL:HB	1:A:270:LEU:HB2	1.99	0.44
1:A:26:GLY:C	1:A:27:TYR:CD1	2.96	0.44
2:E:29:GLN:HA	2:E:61:SER:HB2	1.99	0.44
1:J:22:TYR:CD2	1:J:71:GLU:HG3	2.46	0.44
2:K:41:LYS:HG3	2:K:78:TYR:HE2	1.81	0.44
1:D:8:PHE:HD2	2:E:56:PHE:CE1	2.35	0.44
2:E:21:ASN:HB3	2:E:70:PHE:CE1	2.53	0.44
2:E:34:HIS:ND1	2:E:35:ILE:N	2.66	0.44
1:D:32:GLU:OE2	1:D:35:ARG:NH2	2.45	0.43
1:A:26:GLY:C	1:A:27:TYR:HD1	2.26	0.43
2:E:14:PRO:HA	2:E:15:PRO:HD3	1.80	0.43
2:E:17:ASN:HB3	2:E:74:GLU:OE2	2.17	0.43
1:J:22:TYR:HD2	1:J:71:GLU:HG2	1.83	0.43
2:K:54:MET:HE2	2:K:54:MET:HB2	1.83	0.43
1:J:22:TYR:CD1	1:J:23:ILE:N	2.86	0.43
1:J:81:LEU:HD23	1:J:81:LEU:HA	1.81	0.43
1:J:129:ASP:HB2	1:J:131:LYS:HE3	1.99	0.43
2:K:89:GLU:OE1	2:K:89:GLU:HA	2.19	0.43
2:K:34:HIS:ND1	2:K:34:HIS:C	2.77	0.43
1:A:255:GLN:H	1:A:255:GLN:CD	2.27	0.43
2:H:34:HIS:ND1	2:H:34:HIS:C	2.77	0.43
2:K:97:ARG:H	2:K:97:ARG:HG2	1.65	0.43
2:B:53:ASP:OD2	2:B:53:ASP:N	2.52	0.42
1:J:176:ASN:O	1:J:180:LEU:HB2	2.19	0.42
2:B:21:ASN:HB3	2:B:70:PHE:CE1	2.53	0.42
1:D:47:PRO:HG3	1:D:60:TRP:CZ2	2.54	0.42
1:D:181:ARG:HE	1:D:181:ARG:HB3	1.50	0.42
2:E:53:ASP:OD2	2:E:53:ASP:N	2.52	0.42
1:A:81:LEU:HA	1:A:81:LEU:HD23	1.61	0.42
2:B:29:GLN:HA	2:B:61:SER:HB2	2.00	0.42
1:D:172:LEU:HD12	1:D:172:LEU:HA	1.80	0.42
1:G:255:GLN:H	1:G:255:GLN:CD	2.26	0.42
2:H:53:ASP:OD2	2:H:53:ASP:N	2.52	0.42
2:H:92:THR:HG21	2:H:94:TYR:HE1	1.84	0.42
1:G:47:PRO:HG3	1:G:60:TRP:CZ2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:54:MET:HE2	2:H:54:MET:HB2	1.85	0.42
1:D:274:TRP:O	1:D:276:PRO:HD3	2.20	0.42
2:E:26:TYR:CE2	2:E:28:THR:CG2	3.03	0.42
1:A:181:ARG:HE	1:A:181:ARG:HB3	1.42	0.42
2:B:34:HIS:ND1	2:B:34:HIS:C	2.77	0.41
1:G:213:ILE:O	4:G:301:HOH:O	2.22	0.41
1:G:33:PHE:CD2	1:G:34:VAL:HG23	2.55	0.41
1:J:172:LEU:HD12	1:J:172:LEU:HA	1.79	0.41
1:A:274:TRP:O	1:A:276:PRO:HD3	2.20	0.41
2:E:51:MET:HE3	2:E:51:MET:HB2	2.00	0.41
2:H:14:PRO:HA	2:H:15:PRO:HD3	1.80	0.41
2:K:17:ASN:CB	2:K:74:GLU:OE2	2.69	0.41
2:K:77:THR:C	2:K:78:TYR:HD2	2.29	0.41
2:E:92:THR:HG21	2:E:94:TYR:HE1	1.85	0.41
2:H:92:THR:HG22	2:H:94:TYR:CE1	2.56	0.41
1:G:143:THR:HG23	3:I:9:PHE:HA	2.03	0.41
2:H:94:TYR:N	2:H:94:TYR:CD1	2.88	0.41
2:B:54:MET:HE2	2:B:54:MET:HB2	1.86	0.41
1:G:46:GLU:HA	1:G:47:PRO:HD3	1.95	0.41
1:G:117:ALA:HB2	2:H:60:TRP:CD2	2.56	0.41
1:J:14:ARG:NH2	4:J:309:HOH:O	2.53	0.41
2:K:24:ASN:HB3	2:K:65:LEU:HD11	2.03	0.41
2:H:92:THR:HG22	2:H:94:TYR:HE1	1.86	0.41
1:A:47:PRO:HG3	1:A:60:TRP:CZ2	2.55	0.41
1:A:188:HIS:CE1	1:A:204:TRP:CD1	3.09	0.41
2:E:19:LYS:HA	2:E:20:PRO:HD2	1.94	0.40
1:A:123:TYR:CD1	1:A:123:TYR:C	2.98	0.40
1:D:78:LEU:HD13	1:D:95:LEU:HB2	2.03	0.40
2:H:29:GLN:HA	2:H:61:SER:HB2	2.03	0.40
2:B:19:LYS:HA	2:B:20:PRO:HD2	1.94	0.40
1:G:274:TRP:O	1:G:276:PRO:HD3	2.22	0.40
2:K:77:THR:C	2:K:78:TYR:CD2	3.00	0.40
1:G:98:MET:HE1	2:H:58:LYS:HA	2.02	0.40
2:H:4:THR:HA	2:H:5:PRO:HD3	1.94	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	276/278 (99%)	269 (98%)	7 (2%)	0	100	100
1	D	276/278 (99%)	267 (97%)	9 (3%)	0	100	100
1	G	276/278 (99%)	268 (97%)	8 (3%)	0	100	100
1	J	276/278 (99%)	269 (98%)	7 (2%)	0	100	100
2	B	97/99 (98%)	94 (97%)	3 (3%)	0	100	100
2	E	97/99 (98%)	94 (97%)	3 (3%)	0	100	100
2	H	97/99 (98%)	94 (97%)	3 (3%)	0	100	100
2	K	97/99 (98%)	94 (97%)	3 (3%)	0	100	100
3	C	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
3	F	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
3	I	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
3	L	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
All	All	1520/1544 (98%)	1473 (97%)	47 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	232/232 (100%)	213 (92%)	19 (8%)	10	24
1	D	232/232 (100%)	213 (92%)	19 (8%)	10	24

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	G	232/232 (100%)	214 (92%)	18 (8%)	11	26
1	J	232/232 (100%)	213 (92%)	19 (8%)	10	24
2	B	94/94 (100%)	86 (92%)	8 (8%)	10	22
2	E	94/94 (100%)	87 (93%)	7 (7%)	13	29
2	H	94/94 (100%)	87 (93%)	7 (7%)	13	29
2	K	94/94 (100%)	87 (93%)	7 (7%)	13	29
3	C	9/9 (100%)	9 (100%)	0	100	100
3	F	9/9 (100%)	9 (100%)	0	100	100
3	I	9/9 (100%)	9 (100%)	0	100	100
3	L	9/9 (100%)	8 (89%)	1 (11%)	6	12
All	All	1340/1340 (100%)	1235 (92%)	105 (8%)	11	26

All (105) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	ARG
1	A	41	GLU
1	A	78	LEU
1	A	79	ARG
1	A	105	SER
1	A	109	LEU
1	A	115	GLN
1	A	138	MET
1	A	149	GLN
1	A	170	ARG
1	A	172	LEU
1	A	176	ASN
1	A	178	THR
1	A	181	ARG
1	A	195	SER
1	A	198	GLU
1	A	232	GLU
1	A	258	THR
1	A	272	LEU
2	B	16	GLU
2	B	36	GLU
2	B	45	LYS
2	B	51	MET

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Mol	Chain	Res	Type
2	B	54	MET
2	B	73	THR
2	B	83	LYS
2	B	85	ASP
1	D	21	ARG
1	D	41	GLU
1	D	59	TYR
1	D	78	LEU
1	D	79	ARG
1	D	105	SER
1	D	109	LEU
1	D	115	GLN
1	D	138	MET
1	D	149	GLN
1	D	170	ARG
1	D	172	LEU
1	D	176	ASN
1	D	178	THR
1	D	195	SER
1	D	198	GLU
1	D	232	GLU
1	D	258	THR
1	D	272	LEU
2	E	16	GLU
2	E	45	LYS
2	E	51	MET
2	E	54	MET
2	E	73	THR
2	E	83	LYS
2	E	85	ASP
1	G	21	ARG
1	G	41	GLU
1	G	78	LEU
1	G	79	ARG
1	G	105	SER
1	G	109	LEU
1	G	115	GLN
1	G	138	MET
1	G	170	ARG
1	G	172	LEU
1	G	176	ASN
1	G	178	THR

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Mol	Chain	Res	Type
1	G	195	SER
1	G	198	GLU
1	G	219	LEU
1	G	232	GLU
1	G	258	THR
1	G	272	LEU
2	H	16	GLU
2	H	45	LYS
2	H	51	MET
2	H	54	MET
2	H	73	THR
2	H	83	LYS
2	H	85	ASP
1	J	21	ARG
1	J	41	GLU
1	J	78	LEU
1	J	105	SER
1	J	108	ARG
1	J	109	LEU
1	J	115	GLN
1	J	138	MET
1	J	149	GLN
1	J	170	ARG
1	J	172	LEU
1	J	176	ASN
1	J	178	THR
1	J	195	SER
1	J	198	GLU
1	J	219	LEU
1	J	232	GLU
1	J	258	THR
1	J	272	LEU
2	K	16	GLU
2	K	45	LYS
2	K	51	MET
2	K	54	MET
2	K	73	THR
2	K	83	LYS
2	K	85	ASP
3	L	3	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	96	GLN
1	A	149	GLN
1	A	188	HIS
2	B	13	HIS
1	D	96	GLN
1	D	188	HIS
2	E	13	HIS
1	G	30	ASN
1	G	96	GLN
1	G	188	HIS
2	H	13	HIS
2	H	67	HIS
1	J	96	GLN
1	J	188	HIS
2	K	13	HIS

5.3.3 RNA [i](#)

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	278/278 (100%)	0.26	13 (4%) 36 30	16, 33, 74, 108	10 (3%)
1	D	278/278 (100%)	0.45	21 (7%) 20 16	15, 32, 74, 109	10 (3%)
1	G	278/278 (100%)	0.25	10 (3%) 46 40	17, 33, 74, 108	10 (3%)
1	J	278/278 (100%)	0.31	10 (3%) 46 40	14, 32, 73, 108	10 (3%)
2	B	99/99 (100%)	-0.00	4 (4%) 42 37	14, 30, 64, 89	3 (3%)
2	E	99/99 (100%)	0.05	2 (2%) 65 60	14, 29, 63, 88	3 (3%)
2	H	99/99 (100%)	0.12	4 (4%) 42 37	16, 32, 65, 90	3 (3%)
2	K	99/99 (100%)	0.13	3 (3%) 52 47	15, 31, 65, 93	3 (3%)
3	C	9/9 (100%)	0.37	0 100 100	28, 38, 57, 84	0
3	F	9/9 (100%)	0.42	2 (22%) 2 2	33, 35, 58, 92	0
3	I	9/9 (100%)	0.24	1 (11%) 10 8	21, 34, 62, 83	0
3	L	9/9 (100%)	0.11	0 100 100	22, 30, 37, 66	0
All	All	1544/1544 (100%)	0.26	70 (4%) 38 32	14, 32, 72, 109	52 (3%)

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	181	ARG	5.2
1	J	181	ARG	5.1
1	J	176	ASN	4.8
1	A	54	GLN	4.4
1	D	86	ASN	4.2
2	K	89	GLU	4.0
1	G	54	GLN	3.8
1	D	106	ASP	3.7
1	A	277	PRO	3.6
1	G	277	PRO	3.5
1	J	177	ALA	3.5

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Mol	Chain	Res	Type	RSRZ
1	D	177	ALA	3.3
2	K	48	LYS	3.3
1	J	277	PRO	3.2
2	E	94	TYR	3.1
1	A	107	GLY	3.1
1	A	169	HIS	3.1
1	D	277	PRO	3.1
1	J	170	ARG	3.1
1	D	179	LEU	3.0
1	D	169	HIS	2.9
2	B	48	LYS	2.9
2	E	1	ILE	2.9
1	D	232	GLU	2.9
1	D	180	LEU	2.8
2	B	1	ILE	2.7
1	A	181	ARG	2.7
1	G	105	SER	2.7
1	G	0	ALA	2.7
1	D	170	ARG	2.7
2	B	94	TYR	2.7
1	G	181	ARG	2.7
1	G	169	HIS	2.6
1	D	175	GLY	2.6
2	H	47	PRO	2.6
1	G	88	SER	2.6
1	D	178	THR	2.5
1	J	27	TYR	2.5
1	G	90	GLY	2.5
1	D	54	GLN	2.5
1	D	88	SER	2.4
1	D	105	SER	2.4
1	A	53	GLU	2.4
1	A	59	TYR	2.4
1	J	179	LEU	2.3
1	J	105	SER	2.3
3	F	6	TYR	2.3
1	D	138	MET	2.3
1	A	123	TYR	2.3
1	A	56	GLY	2.2
1	A	65	GLN	2.2
1	D	195	SER	2.2
1	D	27	TYR	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	59	TYR	2.1
2	H	48	LYS	2.1
1	A	180	LEU	2.1
3	F	4	TYR	2.1
1	D	183	ASP	2.1
1	G	178	THR	2.1
1	G	27	TYR	2.1
2	H	94	TYR	2.1
2	B	53	ASP	2.1
1	A	106	ASP	2.0
2	H	89	GLU	2.0
2	K	99	MET	2.0
1	D	194	ARG	2.0
1	A	275	GLU	2.0
1	J	88	SER	2.0
1	J	180	LEU	2.0
3	I	6	TYR	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.