



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 10:12 AM UTC

PDB ID : 8CJ1 / pdb_00008cj1
Title : Urea-based foldamer inhibitor c3u_3 chimera in complex with ASF1 histone chaperone
Authors : Perrin, M.E.; Li, B.; Mbianda, J.; Ropars, V.; Legrand, P.; Douat, C.; Ochsenbein, F.; Guichard, G.
Deposited on : 2023-02-11
Resolution : 2.56 Å(reported)

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

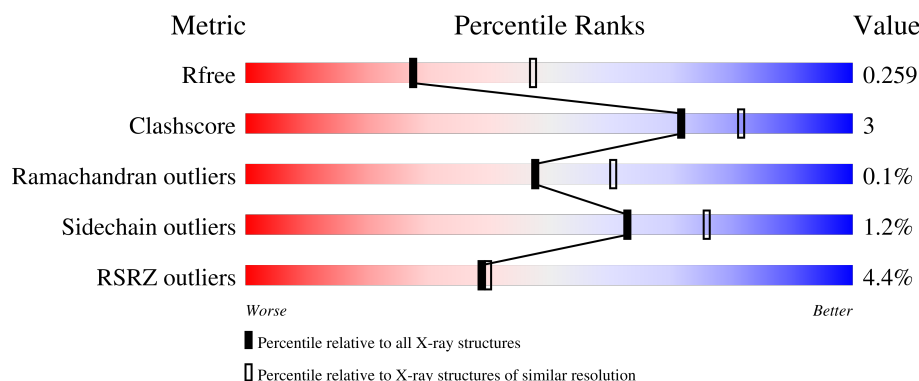
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.56 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	1853 (2.58-2.54)
Clashscore	190562	1897 (2.58-2.54)
Ramachandran outliers	187476	1875 (2.58-2.54)
Sidechain outliers	187428	1875 (2.58-2.54)
RSRZ outliers	180081	1853 (2.58-2.54)

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	158	% 91% 6% .
1	B	158	86% 10% ..
1	C	158	87% 10% ..
1	D	158	% 89% 9% .
1	E	158	3% 85% 12% ..

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Mol	Chain	Length	Quality of chain
1	F	158	6% 90% 8%
1	G	158	6% 91% 6%
1	H	158	15% 86% 11%
2	I	6	67% 17%
2	J	6	17% 50% 33%
2	K	6	33% 67%
2	L	6	33% 67% 17%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10393 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 chaperone ASF1A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	154	Total	C	N	O	S	0	0	0
			1240	796	202	238	4			
1	B	154	Total	C	N	O	S	0	0	0
			1240	796	202	238	4			
1	C	154	Total	C	N	O	S	0	0	0
			1240	796	202	238	4			
1	D	154	Total	C	N	O	S	0	0	0
			1240	796	202	238	4			
1	E	154	Total	C	N	O	S	0	0	0
			1240	796	202	238	4			
1	F	154	Total	C	N	O	S	0	0	0
			1240	796	202	238	4			
1	G	154	Total	C	N	O	S	0	0	0
			1240	796	202	238	4			
1	H	154	Total	C	N	O	S	0	0	0
			1240	796	202	238	4			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP Q9Y294
A	0	ALA	-	expression tag	UNP Q9Y294
B	-1	GLY	-	expression tag	UNP Q9Y294
B	0	ALA	-	expression tag	UNP Q9Y294
C	-1	GLY	-	expression tag	UNP Q9Y294
C	0	ALA	-	expression tag	UNP Q9Y294
D	-1	GLY	-	expression tag	UNP Q9Y294
D	0	ALA	-	expression tag	UNP Q9Y294
E	-1	GLY	-	expression tag	UNP Q9Y294
E	0	ALA	-	expression tag	UNP Q9Y294
F	-1	GLY	-	expression tag	UNP Q9Y294
F	0	ALA	-	expression tag	UNP Q9Y294
G	-1	GLY	-	expression tag	UNP Q9Y294

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Chain	Residue	Modelled	Actual	Comment	Reference
G	0	ALA	-	expression tag	UNP Q9Y294
H	-1	GLY	-	expression tag	UNP Q9Y294
H	0	ALA	-	expression tag	UNP Q9Y294

- Molecule 2 is a protein called c3u_3 chimera inhibitor of histone chaperone ASF1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	I	5	Total	C	N	O	0	0	0
			41	24	12	5			
2	J	5	Total	C	N	O	0	0	0
			41	24	12	5			
2	K	6	Total	C	N	O	0	0	0
			51	31	14	6			
2	L	5	Total	C	N	O	0	0	0
			41	24	12	5			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	53	Total	O	0	0
			53	53		
3	B	48	Total	O	0	0
			48	48		
3	C	42	Total	O	0	0
			42	42		
3	D	39	Total	O	0	0
			39	39		
3	E	24	Total	O	0	0
			24	24		
3	F	34	Total	O	0	0
			34	34		
3	G	33	Total	O	0	0
			33	33		
3	H	22	Total	O	0	0
			22	22		
3	I	1	Total	O	0	0
			1	1		
3	K	1	Total	O	0	0
			1	1		
3	L	2	Total	O	0	0
			2	2		

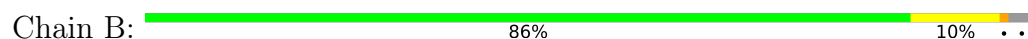
3 Residue-property plots [i](#)

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.

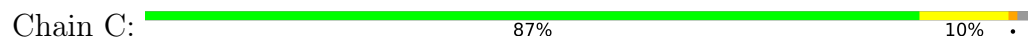
- Molecule 1: Histone chaperone ASF1A



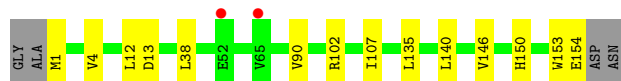
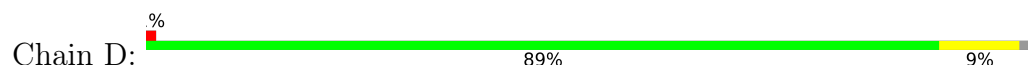
- Molecule 1: Histone chaperone ASF1A



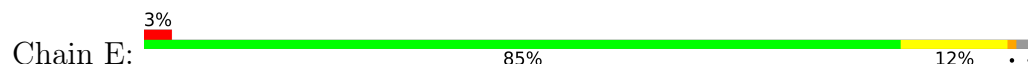
- Molecule 1: Histone chaperone ASF1A



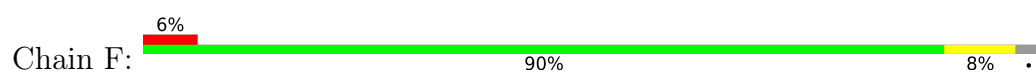
- Molecule 1: Histone chaperone ASF1A



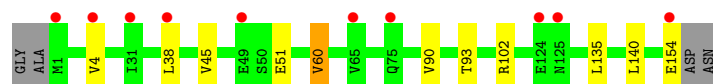
- Molecule 1: Histone chaperone ASF1A



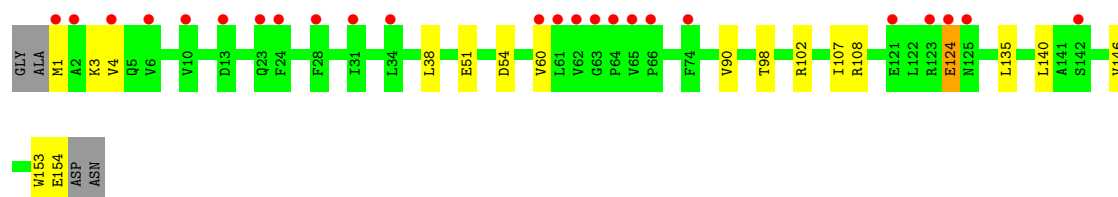
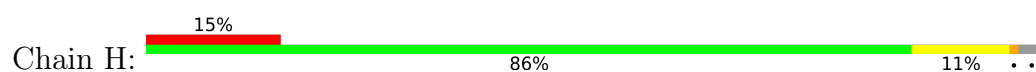
- Molecule 1: Histone chaperone ASF1A



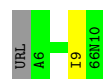
- Molecule 1: Histone chaperone ASF1A



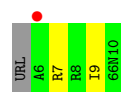
- Molecule 1: Histone chaperone ASF1A



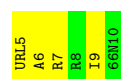
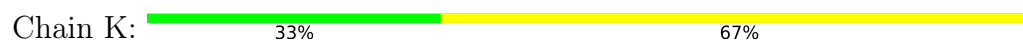
- Molecule 2: c3u_3 chimera inhibitor of histone chaperone ASF1



- Molecule 2: c3u_3 chimera inhibitor of histone chaperone ASF1



- Molecule 2: c3u_3 chimera inhibitor of histone chaperone ASF1



- Molecule 2: c3u_3 chimera inhibitor of histone chaperone ASF1





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	91.56Å 93.22Å 227.68Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	27.88 – 2.56 27.88 – 2.56	Depositor EDS
% Data completeness (in resolution range)	55.7 (27.88-2.56) 58.1 (27.88-2.56)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.06 (at 2.54Å)	Xtriage
Refinement program	BUSTER 2.10.4	Depositor
R, R_{free}	0.225 , 0.280 (Not available) , 0.259	Depositor DCC
R_{free} test set	1782 reflections (2.68%)	wwPDB-VP
Wilson B-factor (Å ²)	40.8	Xtriage
Anisotropy	0.155	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 49.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.054 for k,h,-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	10393	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.75% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 66N, URL

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	0.71	0/1273	0.93	0/1741
1	B	0.73	1/1273 (0.1%)	1.00	3/1741 (0.2%)
1	C	0.79	1/1273 (0.1%)	1.00	4/1741 (0.2%)
1	D	0.71	0/1273	0.98	4/1741 (0.2%)
1	E	0.69	0/1273	0.96	3/1741 (0.2%)
1	F	0.68	0/1273	0.94	2/1741 (0.1%)
1	G	0.69	0/1273	0.95	2/1741 (0.1%)
1	H	0.68	0/1273	1.00	5/1741 (0.3%)
2	I	0.75	0/34	1.46	0/43
2	J	0.60	0/34	1.10	0/43
2	K	0.84	0/34	1.14	0/43
2	L	0.72	0/34	1.06	0/43
All	All	0.71	2/10320 (0.0%)	0.97	23/14100 (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	K	0	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	71	MET	SD-CE	-5.45	1.66	1.79
1	C	71	MET	SD-CE	-5.02	1.67	1.79

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	124	GLU	CB-CG-CD	7.27	124.95	112.60
1	B	35	SER	N-CA-C	-6.49	100.55	109.71
1	H	1	MET	CA-C-N	6.36	130.52	121.42
1	H	1	MET	C-N-CA	6.36	130.52	121.42
1	G	140	LEU	CA-C-N	5.95	128.85	120.28
1	G	140	LEU	C-N-CA	5.95	128.85	120.28
1	D	140	LEU	CA-C-N	5.80	128.64	120.28
1	D	140	LEU	C-N-CA	5.80	128.64	120.28
1	D	153	TRP	CA-C-N	5.44	131.49	121.70
1	D	153	TRP	C-N-CA	5.44	131.49	121.70
1	H	140	LEU	CA-C-N	5.38	128.03	120.28
1	H	140	LEU	C-N-CA	5.38	128.03	120.28
1	E	140	LEU	CA-C-N	5.32	127.94	120.28
1	E	140	LEU	C-N-CA	5.32	127.94	120.28
1	F	140	LEU	CA-C-N	5.31	127.93	120.28
1	F	140	LEU	C-N-CA	5.31	127.93	120.28
1	E	52	GLU	CB-CG-CD	5.25	121.53	112.60
1	B	140	LEU	CA-C-N	5.25	127.84	120.28
1	B	140	LEU	C-N-CA	5.25	127.84	120.28
1	C	140	LEU	CA-C-N	5.23	127.81	120.28
1	C	140	LEU	C-N-CA	5.23	127.81	120.28
1	C	48	ALA	CA-C-N	5.00	127.23	120.38
1	C	48	ALA	C-N-CA	5.00	127.23	120.38

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	K	5	URL	Mainchain,Peptide

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1240	0	1194	8	0
1	B	1240	0	1194	10	0
1	C	1240	0	1194	11	0
1	D	1240	0	1194	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	1240	0	1194	12	0
1	F	1240	0	1194	6	0
1	G	1240	0	1194	9	0
1	H	1240	0	1194	8	0
2	I	41	0	41	2	0
2	J	41	0	41	4	0
2	K	51	0	42	1	0
2	L	41	0	41	1	0
3	A	53	0	0	1	0
3	B	48	0	0	0	0
3	C	42	0	0	0	0
3	D	39	0	0	0	0
3	E	24	0	0	0	0
3	F	34	0	0	0	0
3	G	33	0	0	0	0
3	H	22	0	0	0	0
3	I	1	0	0	0	0
3	K	1	0	0	0	0
3	L	2	0	0	0	0
All	All	10393	0	9717	65	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 (65) 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)
1:G:102:ARG:HH12	1:G:154:GLU:HG3	1.49	0.78
1:B:1:MET:HB2	1:E:51:GLU:OE1	1.84	0.77
1:E:90:VAL:HG22	1:E:135:LEU:HD21	1.69	0.75
1:A:150:HIS:CD2	1:G:51:GLU:HG2	2.26	0.70
1:A:90:VAL:HG22	1:A:135:LEU:HD21	1.74	0.69
1:C:102:ARG:HH12	1:C:154:GLU:HB2	1.61	0.65
1:G:90:VAL:HG22	1:G:135:LEU:HD21	1.78	0.65
1:C:4:VAL:HG21	1:C:38:LEU:HD21	1.80	0.64
1:H:102:ARG:HH12	1:H:154:GLU:HB3	1.62	0.64
1:E:10:VAL:HB	1:E:25:GLU:HB3	1.80	0.63
1:C:122:LEU:HD11	1:C:134:LYS:HD3	1.82	0.61
1:B:90:VAL:HG22	1:B:135:LEU:HD21	1.84	0.59
1:H:3:LYS:HD2	1:H:153:TRP:HB3	1.84	0.58
1:E:102:ARG:HH22	1:E:154:GLU:HB2	1.68	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:150:HIS:ND1	1:H:54:ASP:OD2	2.38	0.56
1:D:90:VAL:HG22	1:D:135:LEU:HD21	1.87	0.56
1:H:98:THR:HG22	1:H:108:ARG:HG3	1.89	0.54
1:F:90:VAL:HG22	1:F:135:LEU:HD21	1.90	0.54
1:E:102:ARG:HH12	1:E:154:GLU:HB3	1.73	0.53
1:H:90:VAL:HG22	1:H:135:LEU:HD21	1.90	0.52
1:A:150:HIS:HD2	1:G:51:GLU:HG2	1.72	0.52
1:F:2:ALA:HB1	1:F:153:TRP:CE2	2.44	0.52
1:C:153:TRP:O	1:C:154:GLU:HB3	2.11	0.51
2:J:7:ARG:HH21	2:J:9:ILE:HD11	1.76	0.50
1:B:1:MET:HE3	1:E:51:GLU:OE1	2.11	0.50
1:C:90:VAL:HG22	1:C:135:LEU:HD21	1.93	0.50
2:K:7:ARG:HH21	2:K:9:ILE:HD11	1.78	0.49
1:D:13:ASP:HA	2:J:9:ILE:HA	1.94	0.49
1:G:45:VAL:O	1:G:93:THR:OG1	2.30	0.48
1:D:4:VAL:HG21	1:D:38:LEU:HD21	1.94	0.48
1:E:30:CYS:SG	1:E:32:GLU:O	2.71	0.48
1:E:9:VAL:HG22	1:E:26:ILE:HG12	1.96	0.48
1:A:40:TRP:HB2	1:A:60:VAL:CG2	2.44	0.47
1:F:50:SER:OG	1:F:52:GLU:HG2	2.15	0.47
1:F:2:ALA:HB1	1:F:153:TRP:CZ2	2.50	0.47
1:C:122:LEU:CD1	1:C:134:LYS:HD3	2.43	0.47
1:C:107:ILE:HD11	1:C:146:VAL:CG1	2.45	0.46
1:D:102:ARG:HH12	1:D:154:GLU:HB3	1.81	0.46
1:B:106:PHE:HB3	1:B:151:ILE:HD13	1.97	0.46
1:A:13:ASP:HA	2:I:9:ILE:HA	1.97	0.45
1:G:4:VAL:HG21	1:G:38:LEU:HD21	1.99	0.45
1:B:107:ILE:HD11	1:B:146:VAL:CG1	2.47	0.45
1:A:111:TYR:CE1	1:A:144:PRO:HB3	2.52	0.44
1:B:28:PHE:CE1	1:B:70:HIS:HB2	2.53	0.44
1:E:107:ILE:HD11	1:E:146:VAL:CG1	2.47	0.44
1:D:107:ILE:HD11	1:D:146:VAL:CG1	2.48	0.43
1:C:1:MET:HE2	1:C:1:MET:HB3	1.86	0.43
1:C:60:VAL:HG11	2:J:9:ILE:HD12	2.00	0.43
1:D:150:HIS:CD2	1:F:54:ASP:OD2	2.71	0.43
1:F:37:ASP:OD1	1:F:63:GLY:HA3	2.18	0.43
1:G:102:ARG:HH22	1:G:154:GLU:H	1.65	0.42
1:B:71:MET:HE3	1:B:71:MET:HB2	1.88	0.42
1:G:60:VAL:HG11	2:L:9:ILE:HD12	2.02	0.42
1:H:4:VAL:HG21	1:H:38:LEU:HD21	2.00	0.42
1:A:40:TRP:HB2	1:A:60:VAL:HG23	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:150:HIS:CD2	1:E:54:ASP:OD2	2.72	0.42
1:D:12:LEU:HB3	2:J:9:ILE:HG23	2.02	0.42
1:B:60:VAL:HG11	2:I:9:ILE:HD12	2.02	0.41
1:G:102:ARG:NH2	1:G:154:GLU:H	2.19	0.41
1:E:102:ARG:HH22	1:E:154:GLU:CB	2.33	0.41
1:E:42:ILE:HG13	1:E:74:PHE:CE2	2.56	0.41
1:B:98:THR:HG22	1:B:108:ARG:HG3	2.02	0.41
1:C:1:MET:HE2	1:H:51:GLU:HB2	2.02	0.41
1:H:107:ILE:HD11	1:H:146:VAL:CG1	2.51	0.41
1:A:46:GLY:HA3	3:A:201:HOH:O	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	152/158 (96%)	149 (98%)	3 (2%)	0	100	100
1	B	152/158 (96%)	149 (98%)	3 (2%)	0	100	100
1	C	152/158 (96%)	150 (99%)	2 (1%)	0	100	100
1	D	152/158 (96%)	151 (99%)	1 (1%)	0	100	100
1	E	152/158 (96%)	149 (98%)	3 (2%)	0	100	100
1	F	152/158 (96%)	149 (98%)	3 (2%)	0	100	100
1	G	152/158 (96%)	149 (98%)	3 (2%)	0	100	100
1	H	152/158 (96%)	149 (98%)	3 (2%)	0	100	100
2	I	2/6 (33%)	2 (100%)	0	0	100	100
2	J	2/6 (33%)	2 (100%)	0	0	100	100
2	K	3/6 (50%)	2 (67%)	0	1 (33%)	0	0
2	L	2/6 (33%)	2 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1225/1288 (95%)	1203 (98%)	21 (2%)	1 (0%)	48 60

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	K	6	ALA

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	139/141 (99%)	138 (99%)	1 (1%)	76 84
1	B	139/141 (99%)	136 (98%)	3 (2%)	45 61
1	C	139/141 (99%)	138 (99%)	1 (1%)	76 84
1	D	139/141 (99%)	138 (99%)	1 (1%)	76 84
1	E	139/141 (99%)	136 (98%)	3 (2%)	45 61
1	F	139/141 (99%)	137 (99%)	2 (1%)	59 73
1	G	139/141 (99%)	138 (99%)	1 (1%)	76 84
1	H	139/141 (99%)	137 (99%)	2 (1%)	59 73
2	I	3/3 (100%)	3 (100%)	0	100 100
2	J	3/3 (100%)	3 (100%)	0	100 100
2	K	3/3 (100%)	3 (100%)	0	100 100
2	L	3/3 (100%)	3 (100%)	0	100 100
All	All	1124/1140 (99%)	1110 (99%)	14 (1%)	63 76

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	CYS
1	B	30	CYS
1	B	52	GLU

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Mol	Chain	Res	Type
1	B	60	VAL
1	C	1	MET
1	D	1	MET
1	E	25	GLU
1	E	98	THR
1	E	150	HIS
1	F	1	MET
1	F	60	VAL
1	G	60	VAL
1	H	60	VAL
1	H	124	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	7	ASN
1	A	143	ASN
1	A	152	ASN
1	B	7	ASN
1	B	23	GLN
1	B	75	GLN
1	B	114	ASN
1	B	150	HIS
1	D	136	GLN
1	D	150	HIS
1	E	150	HIS
1	F	23	GLN
1	F	143	ASN
1	F	150	HIS
1	G	23	GLN
1	G	75	GLN
1	H	7	ASN
1	H	150	HIS

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 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	154/158 (97%)	0.02	2 (1%) 75 76	23, 35, 60, 68	0
1	B	154/158 (97%)	-0.09	0 100 100	22, 36, 56, 64	0
1	C	154/158 (97%)	0.01	0 100 100	18, 37, 59, 70	0
1	D	154/158 (97%)	0.04	2 (1%) 75 76	24, 36, 56, 69	0
1	E	154/158 (97%)	0.48	5 (3%) 50 51	35, 54, 77, 91	0
1	F	154/158 (97%)	0.67	9 (5%) 29 29	28, 51, 71, 78	0
1	G	154/158 (97%)	0.65	10 (6%) 25 25	35, 58, 81, 87	0
1	H	154/158 (97%)	1.05	24 (15%) 5 4	34, 60, 81, 87	0
2	I	4/6 (66%)	0.57	0 100 100	58, 58, 62, 63	0
2	J	4/6 (66%)	1.38	1 (25%) 2 1	63, 65, 71, 73	0
2	K	4/6 (66%)	1.29	0 100 100	78, 84, 86, 87	0
2	L	4/6 (66%)	2.82	2 (50%) 0 0	84, 92, 95, 95	0
All	All	1248/1288 (96%)	0.37	55 (4%) 39 40	18, 46, 75, 95	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	L	6	ALA	6.2
1	H	62	VAL	5.4
1	H	61	LEU	4.5
1	H	64	PRO	4.1
1	A	31	ILE	4.0
1	H	63	GLY	3.9
1	H	2	ALA	3.3
1	F	1	MET	3.2
1	H	28	PHE	3.2
1	H	65	VAL	3.2
1	F	66	PRO	3.1

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Mol	Chain	Res	Type	RSRZ
1	H	4	VAL	2.9
1	H	60	VAL	2.8
1	H	6	VAL	2.8
1	F	103	GLY	2.8
1	H	1	MET	2.8
1	H	142	SER	2.8
1	E	121	GLU	2.7
1	F	65	VAL	2.7
1	H	13	ASP	2.6
1	G	125	ASN	2.6
1	H	125	ASN	2.6
1	H	74	PHE	2.6
1	E	1	MET	2.5
1	D	52	GLU	2.5
1	F	125	ASN	2.5
1	G	4	VAL	2.5
1	H	31	ILE	2.5
2	L	9	ILE	2.5
1	D	65	VAL	2.5
2	J	6	ALA	2.5
1	F	31	ILE	2.4
1	F	2	ALA	2.4
1	G	75	GLN	2.4
1	H	124	GLU	2.3
1	H	34	LEU	2.3
1	E	2	ALA	2.3
1	G	1	MET	2.3
1	G	124	GLU	2.2
1	H	10	VAL	2.2
1	E	125	ASN	2.2
1	F	8	ASN	2.2
1	G	65	VAL	2.2
1	F	28	PHE	2.1
1	G	38	LEU	2.1
1	E	31	ILE	2.1
1	G	49	GLU	2.1
1	H	66	PRO	2.1
1	G	31	ILE	2.1
1	A	101	TYR	2.1
1	H	23	GLN	2.0
1	H	123	ARG	2.0
1	H	24	PHE	2.0

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Mol	Chain	Res	Type	RSRZ
1	G	154	GLU	2.0
1	H	121	GLU	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.