



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 10:55 AM UTC

PDB ID : 3C0D / pdb_00003c0d
Title : Crystal structure of the putative nitrite reductase NADPH (small subunit) oxidoreductase protein Q87HB1. Northeast Structural Genomics Consortium target VpR162
Authors : Kuzin, A.P.; Abashidze, M.; Seetharaman, J.; Vorobiev, S.M.; Wang, D.; Fang, Y.; Owens, L.; Ma, L.-C.; Xiao, R.; Liu, J.; Baran, M.C.; Acton, T.B.; Rost, B.; Montelione, G.T.; Tong, L.; Hunt, J.F.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2008-01-19
Resolution : 2.40 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)

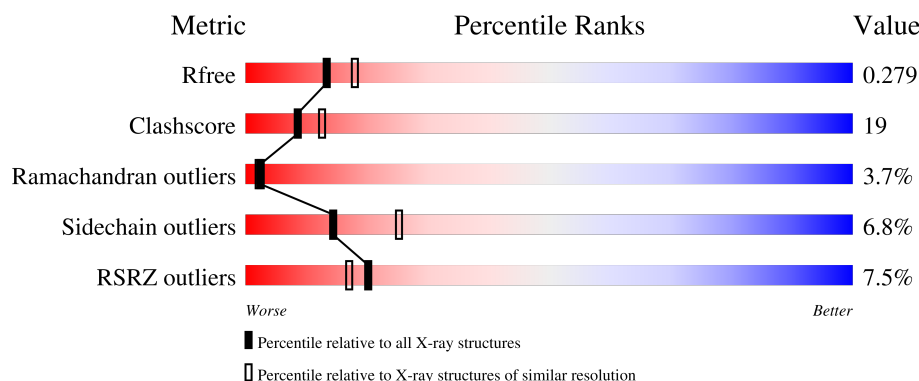
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	119	2% 56% 28% 7% 8%
1	B	119	4% 61% 24% • 10%
1	C	119	6% 53% 29% 5% 11%

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Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

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Mol	Chain	Length	Quality of chain
1	D	119	3% 52% 31% 6% 11%
1	E	119	5% 51% 30% 8% 11%
1	F	119	4% 57% 29% • • 11%
1	G	119	13% 47% 35% 6% • 11%
1	H	119	15% 53% 30% • • 11%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6722 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 Putative nitrite reductase NADPH (Small subunit) oxidoreductase protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	109	Total	C	N	O	S	Se	0	0	0
			852	546	138	160	5	3			
1	B	107	Total	C	N	O	S	Se	0	0	0
			835	535	136	156	5	3			
1	C	106	Total	C	N	O	S	Se	0	0	0
			826	529	134	155	5	3			
1	D	106	Total	C	N	O	S	Se	0	0	0
			826	529	134	155	5	3			
1	E	106	Total	C	N	O	S	Se	0	0	0
			826	529	134	155	5	3			
1	F	106	Total	C	N	O	S	Se	0	0	0
			826	529	134	155	5	3			
1	G	106	Total	C	N	O	S	Se	0	0	0
			826	529	134	155	5	3			
1	H	106	Total	C	N	O	S	Se	0	0	0
			826	529	134	155	5	3			

There are 64 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	112	LEU	-	expression tag	UNP Q87HB1
A	113	GLU	-	expression tag	UNP Q87HB1
A	114	HIS	-	expression tag	UNP Q87HB1
A	115	HIS	-	expression tag	UNP Q87HB1
A	116	HIS	-	expression tag	UNP Q87HB1
A	117	HIS	-	expression tag	UNP Q87HB1
A	118	HIS	-	expression tag	UNP Q87HB1
A	119	HIS	-	expression tag	UNP Q87HB1
B	112	LEU	-	expression tag	UNP Q87HB1
B	113	GLU	-	expression tag	UNP Q87HB1
B	114	HIS	-	expression tag	UNP Q87HB1
B	115	HIS	-	expression tag	UNP Q87HB1

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Chain	Residue	Modelled	Actual	Comment	Reference
B	116	HIS	-	expression tag	UNP Q87HB1
B	117	HIS	-	expression tag	UNP Q87HB1
B	118	HIS	-	expression tag	UNP Q87HB1
B	119	HIS	-	expression tag	UNP Q87HB1
C	112	LEU	-	expression tag	UNP Q87HB1
C	113	GLU	-	expression tag	UNP Q87HB1
C	114	HIS	-	expression tag	UNP Q87HB1
C	115	HIS	-	expression tag	UNP Q87HB1
C	116	HIS	-	expression tag	UNP Q87HB1
C	117	HIS	-	expression tag	UNP Q87HB1
C	118	HIS	-	expression tag	UNP Q87HB1
C	119	HIS	-	expression tag	UNP Q87HB1
D	112	LEU	-	expression tag	UNP Q87HB1
D	113	GLU	-	expression tag	UNP Q87HB1
D	114	HIS	-	expression tag	UNP Q87HB1
D	115	HIS	-	expression tag	UNP Q87HB1
D	116	HIS	-	expression tag	UNP Q87HB1
D	117	HIS	-	expression tag	UNP Q87HB1
D	118	HIS	-	expression tag	UNP Q87HB1
D	119	HIS	-	expression tag	UNP Q87HB1
E	112	LEU	-	expression tag	UNP Q87HB1
E	113	GLU	-	expression tag	UNP Q87HB1
E	114	HIS	-	expression tag	UNP Q87HB1
E	115	HIS	-	expression tag	UNP Q87HB1
E	116	HIS	-	expression tag	UNP Q87HB1
E	117	HIS	-	expression tag	UNP Q87HB1
E	118	HIS	-	expression tag	UNP Q87HB1
E	119	HIS	-	expression tag	UNP Q87HB1
F	112	LEU	-	expression tag	UNP Q87HB1
F	113	GLU	-	expression tag	UNP Q87HB1
F	114	HIS	-	expression tag	UNP Q87HB1
F	115	HIS	-	expression tag	UNP Q87HB1
F	116	HIS	-	expression tag	UNP Q87HB1
F	117	HIS	-	expression tag	UNP Q87HB1
F	118	HIS	-	expression tag	UNP Q87HB1
F	119	HIS	-	expression tag	UNP Q87HB1
G	112	LEU	-	expression tag	UNP Q87HB1
G	113	GLU	-	expression tag	UNP Q87HB1
G	114	HIS	-	expression tag	UNP Q87HB1
G	115	HIS	-	expression tag	UNP Q87HB1
G	116	HIS	-	expression tag	UNP Q87HB1
G	117	HIS	-	expression tag	UNP Q87HB1

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Chain	Residue	Modelled	Actual	Comment	Reference
G	118	HIS	-	expression tag	UNP Q87HB1
G	119	HIS	-	expression tag	UNP Q87HB1
H	112	LEU	-	expression tag	UNP Q87HB1
H	113	GLU	-	expression tag	UNP Q87HB1
H	114	HIS	-	expression tag	UNP Q87HB1
H	115	HIS	-	expression tag	UNP Q87HB1
H	116	HIS	-	expression tag	UNP Q87HB1
H	117	HIS	-	expression tag	UNP Q87HB1
H	118	HIS	-	expression tag	UNP Q87HB1
H	119	HIS	-	expression tag	UNP Q87HB1

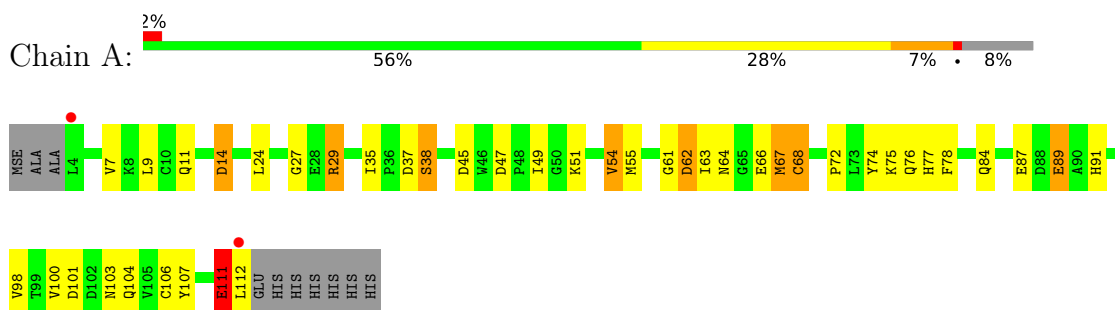
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	9	Total O 9 9	0	0
2	B	13	Total O 13 13	0	0
2	C	6	Total O 6 6	0	0
2	D	9	Total O 9 9	0	0
2	E	21	Total O 21 21	0	0
2	F	9	Total O 9 9	0	0
2	G	10	Total O 10 10	0	0
2	H	2	Total O 2 2	0	0

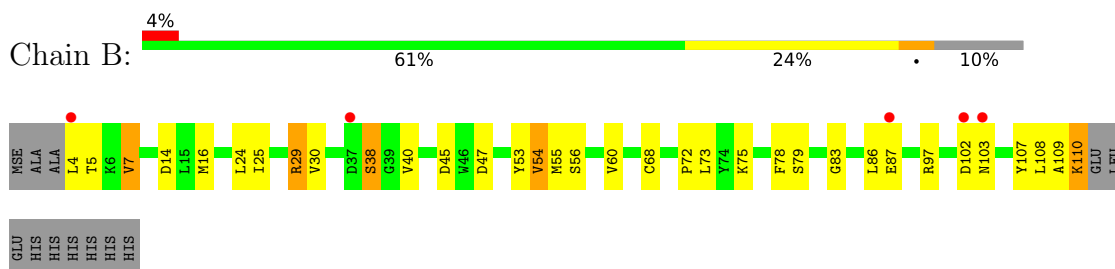
3 Residue-property plots [i](#)

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.

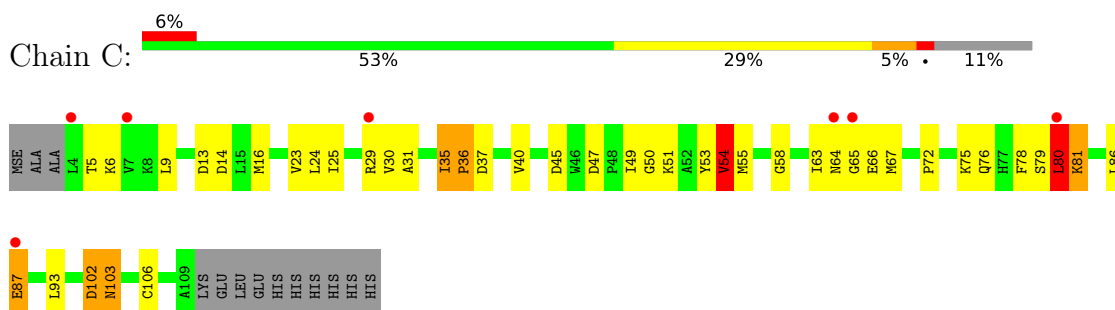
- Molecule 1: Putative nitrite reductase NADPH (Small subunit) oxidoreductase protein



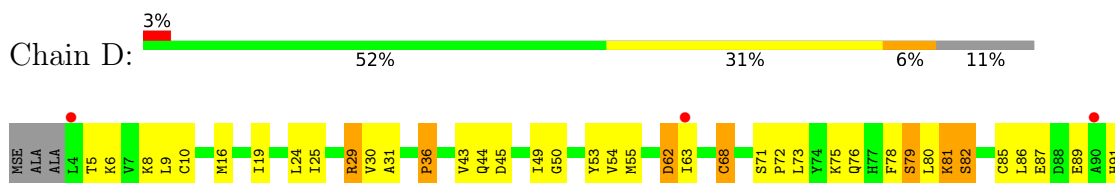
- Molecule 1: Putative nitrite reductase NADPH (Small subunit) oxidoreductase protein



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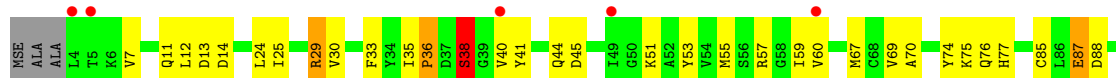




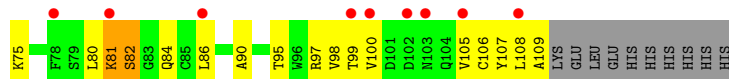
- Molecule 1: Putative nitrite reductase NADPH (Small subunit) oxidoreductase protein



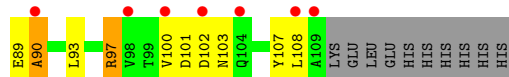
- Molecule 1: Putative nitrite reductase NADPH (Small subunit) oxidoreductase protein



- Molecule 1: Putative nitrite reductase NADPH (Small subunit) oxidoreductase protein



- Molecule 1: Putative nitrite reductase NADPH (Small subunit) oxidoreductase protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	61.67Å 103.46Å 82.95Å 90.00° 91.37° 90.00°	Depositor
Resolution (Å)	19.96 – 2.40 19.96 – 2.40	Depositor EDS
% Data completeness (in resolution range)	73.8 (19.96-2.40) 88.0 (19.96-2.40)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.86 (at 2.39Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.241 , 0.274 0.251 , 0.279	Depositor DCC
R_{free} test set	3791 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å ²)	36.7	Xtriage
Anisotropy	0.960	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 42.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.036 for h,-k,-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	6722	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.35% 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	0.51	0/866	1.05	7/1169 (0.6%)
1	B	0.50	0/849	1.06	6/1146 (0.5%)
1	C	0.49	0/840	1.05	8/1135 (0.7%)
1	D	0.44	0/840	1.02	4/1135 (0.4%)
1	E	0.50	0/840	1.00	4/1135 (0.4%)
1	F	0.43	0/840	0.93	2/1135 (0.2%)
1	G	0.44	0/840	1.01	4/1135 (0.4%)
1	H	0.42	0/840	1.03	4/1135 (0.4%)
All	All	0.47	0/6755	1.02	39/9125 (0.4%)

There are no bond length outliers.

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	89	GLU	N-CA-C	-8.77	101.63	111.71
1	E	14	ASP	N-CA-C	-8.74	102.60	113.18
1	A	14	ASP	N-CA-C	-8.57	102.83	113.28
1	B	14	ASP	N-CA-C	-8.56	102.85	113.38
1	C	103	ASN	N-CA-C	-7.79	100.78	112.54
1	F	75	LYS	N-CA-C	7.56	122.08	112.86
1	C	65	GLY	N-CA-C	-7.46	105.55	115.40
1	A	103	ASN	N-CA-C	-7.39	102.25	112.30
1	G	82	SER	N-CA-C	-6.88	104.88	113.28
1	B	103	ASN	N-CA-C	-6.85	101.27	111.81
1	C	58	GLY	N-CA-C	6.77	122.26	111.64
1	D	75	LYS	N-CA-C	6.73	121.25	113.38
1	H	86	LEU	N-CA-C	-6.69	101.51	110.55
1	B	75	LYS	N-CA-C	6.66	123.85	113.72
1	E	75	LYS	N-CA-C	6.49	122.59	113.56
1	H	103	ASN	N-CA-C	-6.36	104.30	113.21
1	C	35	ILE	CA-C-N	6.23	127.62	119.84
1	C	35	ILE	C-N-CA	6.23	127.62	119.84
1	B	102	ASP	N-CA-C	-6.20	97.60	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	86	LEU	N-CA-C	6.09	117.92	111.28
1	G	14	ASP	N-CA-C	-6.08	106.84	114.56
1	A	51	LYS	N-CA-C	5.91	117.83	110.91
1	H	75	LYS	N-CA-C	5.90	120.06	112.86
1	A	67	MSE	N-CA-C	5.80	119.54	110.32
1	D	68	CYS	N-CA-C	5.79	117.88	109.07
1	G	55	MSE	N-CA-C	5.74	119.80	112.34
1	C	14	ASP	N-CA-C	-5.61	106.49	113.38
1	A	38	SER	N-CA-C	5.59	120.23	113.41
1	E	53	TYR	N-CA-C	-5.41	100.28	108.67
1	B	83	GLY	N-CA-C	-5.39	107.89	115.32
1	B	7	VAL	N-CA-C	5.36	115.61	108.11
1	D	103	ASN	N-CA-C	-5.32	105.76	113.21
1	G	30	VAL	N-CA-C	5.24	116.02	108.48
1	H	102	ASP	N-CA-C	-5.22	105.28	110.97
1	A	75	LYS	N-CA-C	5.22	121.92	110.80
1	A	27	GLY	N-CA-C	5.19	121.25	115.08
1	C	81	LYS	N-CA-C	-5.15	106.94	113.43
1	F	38	SER	N-CA-C	5.07	120.09	113.30
1	C	75	LYS	N-CA-C	5.03	121.52	110.80

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	852	0	848	25	0
1	B	835	0	831	25	0
1	C	826	0	818	28	0
1	D	826	0	818	36	0
1	E	826	0	818	36	0
1	F	826	0	818	33	0
1	G	826	0	818	43	0
1	H	826	0	818	26	0
2	A	9	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	13	0	0	5	0
2	C	6	0	0	0	0
2	D	9	0	0	3	0
2	E	21	0	0	2	0
2	F	9	0	0	1	0
2	G	10	0	0	1	0
2	H	2	0	0	0	0
All	All	6722	0	6587	250	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 19.

All (250) 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:E:29:ARG:NH1	1:E:29:ARG:HB2	1.51	1.23
1:F:29:ARG:HH11	1:F:29:ARG:HB2	1.05	1.17
1:G:29:ARG:NH1	1:G:29:ARG:HB2	1.61	1.14
1:G:29:ARG:HB2	1:G:29:ARG:HH11	1.02	1.11
1:E:29:ARG:HB2	1:E:29:ARG:HH11	0.89	1.03
1:D:29:ARG:HB2	1:D:29:ARG:NH1	1.76	1.00
1:B:29:ARG:HH11	1:B:29:ARG:HB2	1.23	0.99
1:F:29:ARG:HB2	1:F:29:ARG:NH1	1.82	0.93
1:G:81:LYS:HE3	1:G:81:LYS:HA	1.51	0.91
1:G:100:VAL:HG22	1:G:105:VAL:HA	1.53	0.88
1:G:25:ILE:HD12	1:G:30:VAL:HG21	1.56	0.87
1:D:29:ARG:HB2	1:D:29:ARG:HH11	1.33	0.86
1:E:29:ARG:HH11	1:E:29:ARG:CB	1.84	0.84
1:B:29:ARG:HH11	1:B:29:ARG:CB	1.92	0.82
1:G:29:ARG:HD2	1:G:53:TYR:CD1	2.15	0.81
1:E:29:ARG:NH1	1:E:29:ARG:CB	2.41	0.79
1:F:29:ARG:HD2	1:F:53:TYR:CE1	2.16	0.79
1:H:25:ILE:HD12	1:H:30:VAL:HG21	1.66	0.78
1:F:13:ASP:H	1:F:103:ASN:HD21	1.34	0.75
1:C:25:ILE:HD12	1:C:30:VAL:HG21	1.69	0.74
1:E:29:ARG:HG3	2:E:133:HOH:O	1.86	0.73
1:A:61:GLY:HA2	1:C:16:MSE:HE1	1.70	0.72
1:C:55:MSE:HE2	1:C:93:LEU:HD11	1.72	0.72
1:G:16:MSE:HE3	1:G:17:PRO:HD3	1.71	0.72
1:G:29:ARG:HH11	1:G:29:ARG:CB	1.93	0.71
1:G:9:LEU:HA	2:G:123:HOH:O	1.90	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:97:ARG:HH11	1:F:97:ARG:HG2	1.56	0.69
1:D:68:CYS:HB3	1:D:79:SER:HA	1.74	0.69
1:B:29:ARG:HD2	1:B:53:TYR:CD1	2.28	0.69
1:D:68:CYS:HB2	1:D:78:PHE:O	1.92	0.69
1:G:59:ILE:H	1:G:59:ILE:HD12	1.58	0.68
1:F:38:SER:HB2	1:F:41:TYR:OH	1.95	0.67
1:E:29:ARG:HD3	2:E:126:HOH:O	1.94	0.67
1:E:29:ARG:HD2	1:E:53:TYR:CD1	2.30	0.67
1:B:29:ARG:HD2	1:B:53:TYR:CE1	2.32	0.65
1:H:79:SER:OG	1:H:82:SER:HB3	1.97	0.65
1:A:54:VAL:HG22	1:A:72:PRO:HG3	1.78	0.64
1:F:74:TYR:HB3	1:F:76:GLN:HE22	1.60	0.64
1:B:29:ARG:HD3	2:B:120:HOH:O	1.98	0.64
1:F:74:TYR:CB	1:F:76:GLN:HE22	2.11	0.64
1:B:4:LEU:HB2	1:B:110:LYS:HD2	1.80	0.64
1:C:45:ASP:HA	1:C:55:MSE:HB2	1.80	0.63
1:D:62:ASP:C	1:D:63:ILE:HD12	2.24	0.63
1:D:62:ASP:O	1:D:63:ILE:HD12	1.99	0.62
1:A:111:GLU:O	1:A:112:LEU:HB2	1.98	0.62
1:B:7:VAL:HB	1:B:107:TYR:CE1	2.35	0.61
1:B:54:VAL:HG13	1:B:72:PRO:HD2	1.81	0.61
1:G:43:VAL:HG12	1:G:95:THR:HG22	1.81	0.61
1:D:29:ARG:HD2	1:D:53:TYR:CD1	2.36	0.61
1:C:54:VAL:HG13	1:C:72:PRO:HD2	1.80	0.61
1:G:59:ILE:HD12	1:G:59:ILE:N	2.15	0.61
1:G:54:VAL:HG22	1:G:72:PRO:HG2	1.83	0.61
1:C:64:ASN:C	1:C:66:GLU:H	2.09	0.60
1:D:43:VAL:HG11	1:D:93:LEU:HD13	1.83	0.60
1:A:68:CYS:HB3	1:A:78:PHE:C	2.27	0.60
1:D:25:ILE:HD12	1:D:30:VAL:HG21	1.83	0.60
1:G:29:ARG:HD2	1:G:53:TYR:CE1	2.36	0.60
1:B:68:CYS:HB3	1:B:79:SER:HA	1.82	0.60
1:F:29:ARG:HH11	1:F:29:ARG:CB	1.97	0.60
1:F:12:LEU:HD22	1:F:100:VAL:HG11	1.84	0.59
1:C:5:THR:HG22	1:C:6:LYS:N	2.18	0.59
1:F:57:ARG:HA	1:F:57:ARG:HH11	1.67	0.58
1:H:77:HIS:HB2	1:H:86:LEU:HB2	1.85	0.58
1:C:86:LEU:O	1:C:87:GLU:HB2	2.03	0.58
1:B:5:THR:HB	1:B:109:ALA:HB3	1.84	0.58
1:H:5:THR:O	1:H:108:LEU:HA	2.05	0.57
1:B:87:GLU:HB2	2:B:130:HOH:O	2.02	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:63:ILE:HG21	1:C:86:LEU:HD23	1.87	0.57
1:G:57:ARG:HH11	1:G:57:ARG:HA	1.68	0.57
1:A:77:HIS:HD2	1:A:87:GLU:OE2	1.88	0.57
1:D:29:ARG:HD2	1:D:53:TYR:CE1	2.40	0.56
1:F:97:ARG:HG2	1:F:97:ARG:NH1	2.19	0.56
1:B:68:CYS:HB2	1:B:78:PHE:O	2.05	0.56
1:G:37:ASP:CG	1:G:38:SER:H	2.14	0.56
1:F:11:GLN:HB2	1:F:14:ASP:OD2	2.06	0.56
1:F:25:ILE:HD12	1:F:30:VAL:HG21	1.87	0.56
1:E:9:LEU:HD13	1:E:107:TYR:HD2	1.71	0.56
1:C:35:ILE:HG12	1:C:67:MSE:HG2	1.88	0.56
1:D:55:MSE:HE2	1:D:93:LEU:HD11	1.88	0.55
1:D:49:ILE:HD13	1:D:76:GLN:HG2	1.87	0.55
1:C:49:ILE:HG21	1:C:76:GLN:NE2	2.23	0.54
1:C:67:MSE:HB3	1:C:80:LEU:HD21	1.88	0.54
1:D:43:VAL:CG1	1:D:93:LEU:HD13	2.37	0.54
1:D:54:VAL:HG22	1:D:54:VAL:O	2.06	0.54
1:A:98:VAL:HG13	1:A:106:CYS:O	2.07	0.54
1:D:29:ARG:HD3	2:D:120:HOH:O	2.06	0.54
1:D:81:LYS:NZ	1:D:97:ARG:HH11	2.05	0.54
1:E:97:ARG:HE	1:E:97:ARG:HA	1.72	0.54
1:F:44:GLN:HG2	1:F:93:LEU:HB3	1.90	0.54
1:F:29:ARG:HD2	1:F:53:TYR:CD1	2.42	0.54
1:H:86:LEU:O	1:H:87:GLU:C	2.51	0.54
1:H:55:MSE:HE2	1:H:93:LEU:HD11	1.88	0.53
1:C:78:PHE:CD2	1:C:93:LEU:HG	2.42	0.53
1:B:29:ARG:HH11	1:B:29:ARG:CG	2.21	0.53
1:G:4:LEU:HD23	1:G:5:THR:N	2.22	0.53
1:B:54:VAL:HG11	1:B:73:LEU:HG	1.90	0.53
1:D:30:VAL:HG12	1:D:31:ALA:N	2.24	0.53
1:A:68:CYS:HB3	1:A:78:PHE:O	2.09	0.52
1:G:11:GLN:O	1:G:14:ASP:HB2	2.09	0.52
1:C:23:VAL:HG23	1:C:30:VAL:HB	1.91	0.52
1:E:97:ARG:HA	1:E:97:ARG:NE	2.24	0.52
1:D:79:SER:OG	1:D:82:SER:HB3	2.09	0.52
1:E:9:LEU:HD13	1:E:107:TYR:CD2	2.44	0.52
1:E:63:ILE:HG22	1:E:64:ASN:OD1	2.10	0.52
1:D:54:VAL:HG22	1:D:72:PRO:HG2	1.92	0.52
1:B:25:ILE:HD12	1:B:30:VAL:HG21	1.90	0.52
1:D:29:ARG:CG	2:D:120:HOH:O	2.58	0.52
1:G:9:LEU:HD13	1:G:106:CYS:HA	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:45:ASP:HA	1:D:55:MSE:HB2	1.92	0.52
1:E:29:ARG:HD2	1:E:53:TYR:CE1	2.44	0.51
1:G:29:ARG:NH1	1:G:29:ARG:CB	2.55	0.51
1:A:61:GLY:O	1:A:67:MSE:HB2	2.11	0.51
1:F:51:LYS:HE2	2:F:134:HOH:O	2.10	0.51
1:G:7:VAL:HG23	1:G:109:ALA:HB2	1.92	0.51
1:F:12:LEU:HB3	1:F:103:ASN:HD22	1.76	0.51
1:H:17:PRO:O	1:H:18:PHE:HB2	2.09	0.51
1:A:54:VAL:HG22	1:A:72:PRO:CG	2.39	0.51
1:D:5:THR:O	1:D:108:LEU:HA	2.10	0.51
1:E:11:GLN:O	1:E:14:ASP:HB2	2.11	0.51
1:B:60:VAL:HG22	2:B:128:HOH:O	2.10	0.51
1:F:33:PHE:O	1:F:40:VAL:HA	2.11	0.51
1:H:85:CYS:SG	1:H:86:LEU:N	2.83	0.50
1:E:84:GLN:NE2	1:E:89:GLU:HA	2.26	0.50
1:D:81:LYS:HZ2	1:D:97:ARG:HH11	1.58	0.50
1:F:29:ARG:HD2	1:F:53:TYR:CZ	2.47	0.50
1:G:67:MSE:HG3	1:G:80:LEU:HD12	1.94	0.50
1:D:29:ARG:HB2	1:D:29:ARG:CZ	2.40	0.50
1:E:78:PHE:CD2	1:E:93:LEU:HG	2.47	0.50
1:H:38:SER:HB3	1:H:41:TYR:OH	2.12	0.50
1:A:9:LEU:HD13	1:A:107:TYR:HD2	1.78	0.49
1:B:86:LEU:N	1:B:86:LEU:HD12	2.27	0.49
1:E:81:LYS:NZ	1:E:97:ARG:HE	2.11	0.49
1:B:5:THR:O	1:B:108:LEU:HA	2.13	0.49
1:H:29:ARG:HG2	1:H:53:TYR:CD1	2.48	0.49
1:C:5:THR:HG22	1:C:6:LYS:H	1.76	0.49
1:F:70:ALA:HB2	1:F:77:HIS:CE1	2.48	0.49
1:D:43:VAL:HG12	1:D:44:GLN:O	2.13	0.48
1:A:62:ASP:OD1	1:A:62:ASP:N	2.45	0.48
1:E:78:PHE:CE2	1:E:85:CYS:HB2	2.48	0.48
1:F:69:VAL:HG13	1:F:69:VAL:O	2.12	0.48
1:C:64:ASN:C	1:C:66:GLU:N	2.71	0.48
1:F:29:ARG:NH1	1:F:29:ARG:CB	2.66	0.48
1:H:86:LEU:O	1:H:88:ASP:N	2.46	0.48
1:A:45:ASP:OD2	2:A:127:HOH:O	2.20	0.48
1:A:84:GLN:NE2	1:A:89:GLU:HA	2.29	0.47
1:G:86:LEU:N	1:G:86:LEU:HD12	2.29	0.47
1:D:16:MSE:HB3	1:D:19:ILE:HD12	1.97	0.47
1:B:29:ARG:CG	1:B:29:ARG:NH1	2.76	0.47
1:E:12:LEU:HD21	1:E:34:TYR:CZ	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:82:SER:C	1:G:84:GLN:H	2.23	0.47
1:B:38:SER:HB2	2:B:132:HOH:O	2.14	0.47
1:D:5:THR:HG22	1:D:6:LYS:N	2.30	0.47
1:D:9:LEU:O	1:D:10:CYS:HB3	2.14	0.47
1:G:18:PHE:HB3	1:G:60:VAL:HG21	1.96	0.47
1:A:63:ILE:HG22	1:A:64:ASN:OD1	2.15	0.46
1:C:29:ARG:HG3	1:C:53:TYR:CE1	2.50	0.46
1:A:7:VAL:HB	1:A:107:TYR:CE2	2.50	0.46
1:G:18:PHE:HB3	1:G:60:VAL:CG2	2.46	0.46
1:C:37:ASP:OD2	1:C:37:ASP:N	2.49	0.46
1:H:84:GLN:NE2	1:H:85:CYS:N	2.64	0.46
1:F:45:ASP:HA	1:F:55:MSE:HB2	1.97	0.46
1:A:74:TYR:HB3	1:A:76:GLN:HE22	1.81	0.46
1:C:30:VAL:HG12	1:C:31:ALA:N	2.31	0.46
1:E:50:GLY:O	1:E:51:LYS:HB2	2.16	0.46
1:F:59:ILE:N	1:F:59:ILE:HD12	2.31	0.46
1:G:47:ASP:CG	1:G:49:ILE:O	2.59	0.46
1:A:11:GLN:O	1:A:14:ASP:HB2	2.16	0.45
1:D:85:CYS:O	1:D:86:LEU:HD12	2.17	0.45
1:H:17:PRO:HG3	1:H:34:TYR:CD2	2.51	0.45
1:G:97:ARG:HD2	1:G:97:ARG:HA	1.80	0.45
1:G:98:VAL:HG12	1:G:99:THR:N	2.32	0.45
1:E:15:LEU:O	1:E:16:MSE:HE2	2.16	0.45
1:G:41:TYR:CD2	1:G:81:LYS:HD2	2.52	0.45
1:H:12:LEU:HD11	1:H:34:TYR:CE1	2.51	0.45
1:G:9:LEU:HD11	1:G:107:TYR:CD1	2.52	0.45
1:E:9:LEU:O	1:E:10:CYS:HB3	2.16	0.44
1:G:4:LEU:HD21	1:G:108:LEU:HD22	1.99	0.44
1:H:100:VAL:HG12	1:H:101:ASP:N	2.31	0.44
1:F:59:ILE:HD12	1:F:59:ILE:H	1.83	0.44
1:H:29:ARG:HD3	1:H:29:ARG:HA	1.79	0.44
1:H:81:LYS:O	1:H:81:LYS:HD3	2.18	0.44
1:B:47:ASP:C	1:B:47:ASP:OD1	2.60	0.44
1:C:50:GLY:O	1:C:51:LYS:HB2	2.17	0.44
1:H:88:ASP:OD1	1:H:90:ALA:HB3	2.18	0.44
1:A:35:ILE:HB	1:A:38:SER:HB2	1.99	0.44
1:E:41:TYR:CD1	1:E:81:LYS:HB2	2.53	0.44
1:F:13:ASP:H	1:F:103:ASN:ND2	2.09	0.44
1:G:9:LEU:HD11	1:G:107:TYR:HD1	1.82	0.44
1:E:47:ASP:C	1:E:47:ASP:OD1	2.59	0.44
1:B:97:ARG:HD2	1:B:97:ARG:HA	1.74	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:9:LEU:HB2	1:C:106:CYS:HA	2.00	0.43
1:D:78:PHE:CD1	1:D:93:LEU:HG	2.53	0.43
1:A:9:LEU:HD13	1:A:107:TYR:CD2	2.52	0.43
1:A:100:VAL:HA	1:A:104:GLN:O	2.18	0.43
1:C:55:MSE:HE2	1:C:93:LEU:CD1	2.43	0.43
1:D:50:GLY:HA3	1:D:73:LEU:HD13	2.01	0.43
1:A:47:ASP:OD1	1:A:47:ASP:C	2.61	0.43
1:E:43:VAL:HB	1:E:93:LEU:HD13	1.99	0.43
1:A:49:ILE:HD11	1:A:91:HIS:CD2	2.54	0.43
1:A:54:VAL:HG22	1:A:54:VAL:O	2.18	0.43
1:C:25:ILE:CD1	1:C:30:VAL:HG21	2.44	0.43
1:C:79:SER:C	1:C:81:LYS:H	2.27	0.43
1:D:25:ILE:CD1	1:D:30:VAL:HG21	2.47	0.43
1:E:17:PRO:O	1:E:18:PHE:HB2	2.18	0.43
1:F:35:ILE:HD12	1:F:67:MSE:HE2	2.01	0.43
1:E:55:MSE:HE2	1:E:93:LEU:HD11	2.01	0.43
1:G:29:ARG:HG2	1:G:45:ASP:HB3	2.01	0.43
1:D:71:SER:HA	1:D:72:PRO:HD3	1.87	0.43
1:E:30:VAL:HG22	1:E:96:TRP:CZ3	2.53	0.42
1:E:71:SER:O	1:E:75:LYS:HA	2.18	0.42
1:F:35:ILE:CD1	1:F:67:MSE:HE2	2.48	0.42
1:F:85:CYS:HB3	1:F:88:ASP:O	2.20	0.42
1:E:5:THR:HG22	1:E:6:LYS:N	2.33	0.42
1:E:40:VAL:O	1:E:40:VAL:HG13	2.18	0.42
1:F:87:GLU:OE1	1:F:87:GLU:N	2.51	0.42
1:H:97:ARG:O	1:H:107:TYR:HA	2.20	0.42
1:C:47:ASP:CG	1:C:50:GLY:H	2.27	0.42
1:E:29:ARG:HD2	1:E:53:TYR:CG	2.55	0.42
1:E:63:ILE:HG22	1:E:64:ASN:N	2.34	0.42
1:E:81:LYS:O	1:E:82:SER:CB	2.68	0.42
1:E:98:VAL:HG12	1:E:99:THR:N	2.35	0.42
1:H:17:PRO:C	1:H:19:ILE:H	2.28	0.42
1:B:45:ASP:HA	1:B:55:MSE:HB2	2.02	0.42
1:D:68:CYS:HA	1:D:80:LEU:HG	2.01	0.42
1:G:4:LEU:HD21	1:G:108:LEU:HB3	2.01	0.42
1:G:55:MSE:HE1	1:G:71:SER:CA	2.49	0.42
1:G:63:ILE:C	1:G:65:GLY:H	2.28	0.42
1:B:54:VAL:HG21	1:G:73:LEU:HD21	2.02	0.41
1:D:49:ILE:CD1	1:D:76:GLN:HG2	2.50	0.41
1:F:60:VAL:HG12	1:F:69:VAL:HB	2.02	0.41
1:H:47:ASP:CG	1:H:49:ILE:O	2.64	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:63:ILE:HG22	1:G:64:ASN:N	2.35	0.41
1:C:102:ASP:HB3	1:C:103:ASN:H	1.69	0.41
1:C:45:ASP:O	1:C:55:MSE:HG2	2.21	0.41
1:D:8:LYS:HB3	2:D:123:HOH:O	2.20	0.41
1:C:54:VAL:HG13	1:C:72:PRO:CD	2.49	0.41
1:G:63:ILE:HD12	1:G:86:LEU:CD2	2.50	0.41
1:A:29:ARG:CA	1:A:29:ARG:HE	2.29	0.41
1:G:71:SER:O	1:G:75:LYS:HG2	2.20	0.41
1:H:41:TYR:CD1	1:H:81:LYS:HB2	2.56	0.41
1:H:86:LEU:C	1:H:86:LEU:HD23	2.46	0.41
1:B:87:GLU:N	2:B:130:HOH:O	2.53	0.41
1:G:37:ASP:CG	1:G:38:SER:N	2.76	0.40
1:H:54:VAL:O	1:H:54:VAL:HG22	2.21	0.40
1:A:45:ASP:HA	1:A:55:MSE:HB2	2.02	0.40
1:H:77:HIS:O	1:H:86:LEU:N	2.54	0.40
1:F:7:VAL:HB	1:F:107:TYR:CE2	2.57	0.40
1:G:7:VAL:HB	1:G:107:TYR:CE1	2.56	0.40
1:E:29:ARG:HB2	1:E:29:ARG:CZ	2.37	0.40
1:H:55:MSE:C	1:H:57:ARG:N	2.80	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	107/119 (90%)	90 (84%)	13 (12%)	4 (4%)	2	2
1	B	105/119 (88%)	93 (89%)	11 (10%)	1 (1%)	12	20
1	C	104/119 (87%)	88 (85%)	11 (11%)	5 (5%)	2	1
1	D	104/119 (87%)	87 (84%)	13 (12%)	4 (4%)	2	2
1	E	104/119 (87%)	92 (88%)	8 (8%)	4 (4%)	2	2

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	104/119 (87%)	89 (86%)	14 (14%)	1 (1%)	12	20
1	G	104/119 (87%)	84 (81%)	16 (15%)	4 (4%)	2	2
1	H	104/119 (87%)	83 (80%)	13 (12%)	8 (8%)	1	0
All	All	836/952 (88%)	706 (84%)	99 (12%)	31 (4%)	2	2

All (31) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	87	GLU
1	D	87	GLU
1	E	82	SER
1	H	63	ILE
1	H	87	GLU
1	H	88	ASP
1	A	111	GLU
1	C	102	ASP
1	D	82	SER
1	A	89	GLU
1	D	36	PRO
1	D	91	HIS
1	G	51	LYS
1	G	73	LEU
1	H	82	SER
1	H	90	ALA
1	B	38	SER
1	C	36	PRO
1	C	54	VAL
1	C	80	LEU
1	E	54	VAL
1	F	36	PRO
1	G	38	SER
1	A	54	VAL
1	A	101	ASP
1	E	36	PRO
1	E	51	LYS
1	G	90	ALA
1	H	76	GLN
1	H	65	GLY
1	H	36	PRO

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	94/98 (96%)	87 (93%)	7 (7%)	13	22
1	B	92/98 (94%)	85 (92%)	7 (8%)	12	21
1	C	91/98 (93%)	85 (93%)	6 (7%)	15	26
1	D	91/98 (93%)	84 (92%)	7 (8%)	12	20
1	E	91/98 (93%)	83 (91%)	8 (9%)	9	15
1	F	91/98 (93%)	86 (94%)	5 (6%)	19	34
1	G	91/98 (93%)	88 (97%)	3 (3%)	33	55
1	H	91/98 (93%)	84 (92%)	7 (8%)	12	20
All	All	732/784 (93%)	682 (93%)	50 (7%)	14	25

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	LEU
1	A	29	ARG
1	A	37	ASP
1	A	62	ASP
1	A	66	GLU
1	A	68	CYS
1	A	111	GLU
1	B	16	MSE
1	B	24	LEU
1	B	29	ARG
1	B	40	VAL
1	B	54	VAL
1	B	56	SER
1	B	110	LYS
1	C	13	ASP
1	C	24	LEU
1	C	36	PRO
1	C	40	VAL
1	C	54	VAL

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Mol	Chain	Res	Type
1	C	80	LEU
1	D	24	LEU
1	D	29	ARG
1	D	36	PRO
1	D	62	ASP
1	D	79	SER
1	D	81	LYS
1	D	98	VAL
1	E	24	LEU
1	E	29	ARG
1	E	36	PRO
1	E	49	ILE
1	E	81	LYS
1	E	87	GLU
1	E	97	ARG
1	E	101	ASP
1	F	24	LEU
1	F	29	ARG
1	F	36	PRO
1	F	38	SER
1	F	87	GLU
1	G	29	ARG
1	G	55	MSE
1	G	81	LYS
1	H	13	ASP
1	H	28	GLU
1	H	81	LYS
1	H	86	LEU
1	H	87	GLU
1	H	89	GLU
1	H	97	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	76	GLN
1	A	77	HIS
1	B	11	GLN
1	B	91	HIS
1	C	11	GLN
1	C	77	HIS
1	C	84	GLN

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Mol	Chain	Res	Type
1	D	11	GLN
1	D	91	HIS
1	E	103	ASN
1	F	64	ASN
1	F	76	GLN
1	F	77	HIS
1	F	84	GLN
1	F	103	ASN
1	G	104	GLN
1	H	91	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	106/119 (89%)	0.26	2 (1%) 66 62	24, 44, 71, 83	0
1	B	104/119 (87%)	0.36	5 (4%) 35 32	25, 47, 73, 77	0
1	C	103/119 (86%)	0.62	7 (6%) 23 20	23, 52, 81, 86	1 (0%)
1	D	103/119 (86%)	0.43	3 (2%) 53 50	28, 50, 78, 86	0
1	E	103/119 (86%)	0.29	6 (5%) 29 25	26, 44, 66, 77	0
1	F	103/119 (86%)	0.67	5 (4%) 35 31	32, 59, 85, 92	0
1	G	103/119 (86%)	0.99	16 (15%) 5 4	28, 68, 92, 97	0
1	H	103/119 (86%)	1.18	18 (17%) 4 3	35, 76, 93, 104	0
All	All	828/952 (86%)	0.60	62 (7%) 20 17	23, 52, 85, 104	1 (0%)

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	63	ILE	4.0
1	D	4	LEU	3.9
1	F	4	LEU	3.8
1	H	90	ALA	3.8
1	A	112	LEU	3.7
1	D	90	ALA	3.3
1	C	4	LEU	3.2
1	H	65	GLY	3.1
1	H	100	VAL	3.0
1	H	108	LEU	3.0
1	C	80	LEU	2.9
1	E	4	LEU	2.9
1	C	29	ARG	2.9
1	H	87	GLU	2.9
1	F	49	ILE	2.8
1	G	13	ASP	2.8

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Mol	Chain	Res	Type	RSRZ
1	H	81	LYS	2.8
1	G	14	ASP	2.8
1	A	4	LEU	2.7
1	B	87	GLU	2.7
1	E	101	ASP	2.7
1	H	69	VAL	2.7
1	H	109	ALA	2.7
1	B	37	ASP	2.7
1	G	81	LYS	2.6
1	H	86	LEU	2.6
1	G	100	VAL	2.6
1	F	5	THR	2.6
1	E	81	LYS	2.6
1	G	5	THR	2.5
1	G	99	THR	2.5
1	B	4	LEU	2.5
1	G	78	PHE	2.5
1	H	63	ILE	2.4
1	D	63	ILE	2.4
1	G	108	LEU	2.4
1	C	7	VAL	2.4
1	C	64	ASN	2.4
1	G	103	ASN	2.4
1	B	103	ASN	2.4
1	H	4	LEU	2.3
1	H	34	TYR	2.3
1	H	104	GLN	2.3
1	G	68	CYS	2.3
1	H	12	LEU	2.2
1	G	102	ASP	2.2
1	H	102	ASP	2.2
1	F	40	VAL	2.1
1	H	98	VAL	2.1
1	E	64	ASN	2.1
1	G	86	LEU	2.1
1	C	87	GLU	2.1
1	E	37	ASP	2.1
1	C	65	GLY	2.1
1	E	63	ILE	2.1
1	F	60	VAL	2.1
1	H	15	LEU	2.1
1	G	39	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	G	105	VAL	2.0
1	G	4	LEU	2.0
1	H	66	GLU	2.0
1	B	102	ASP	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.