



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:57 AM UTC

PDB ID : 1GZ0 / pdb_00001gz0
Title : 23S RIBOSOMAL RNA G2251 2'O-METHYLTRANSFERASE RLMB
Authors : Michel, G.; Cygler, M.
Deposited on : 2002-05-03
Resolution : 2.50 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

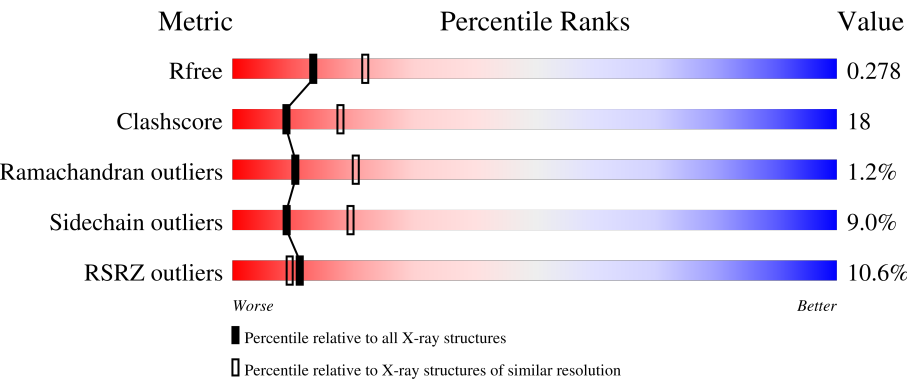
MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry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

1 Overall quality at a glance 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}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	253	13% 60% 31% •••
1	B	253	13% 56% 34% 6% •
1	C	253	6% 62% 28% 5% •
1	D	253	8% 60% 32% ••
1	E	253	42% 20% 34% •

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Mol	Chain	Length	Quality of chain
1	F	253	22% 62% 30% 7%
1	G	253	45% 17% • 34%
1	H	253	10% 71% 26% ••

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 13953 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 HYPOTHETICAL TRNA/RRNA METHYLTRANSFERASE YJFH.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	242	Total	C	N	O	S	Se	4	0	0
			1851	1153	339	348	4	7			
1	B	244	Total	C	N	O	S	Se	0	0	0
			1869	1164	343	350	4	8			
1	C	242	Total	C	N	O	S	Se	0	0	0
			1851	1153	339	348	4	7			
1	D	244	Total	C	N	O	S	Se	0	0	0
			1869	1164	343	350	4	8			
1	E	166	Total	C	N	O	S	Se	0	0	0
			1248	770	228	240	4	6			
1	F	253	Total	C	N	O	S	Se	0	0	0
			1928	1199	355	362	4	8			
1	G	166	Total	C	N	O	S	Se	0	0	0
			1248	770	228	240	4	6			
1	H	250	Total	C	N	O	S	Se	0	0	0
			1912	1191	352	357	4	8			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	14	Total	O	0	0
			14	14		
2	B	19	Total	O	0	0
			19	19		
2	C	14	Total	O	0	0
			14	14		
2	D	24	Total	O	0	0
			24	24		
2	E	28	Total	O	0	0
			28	28		
2	F	38	Total	O	0	0
			38	38		

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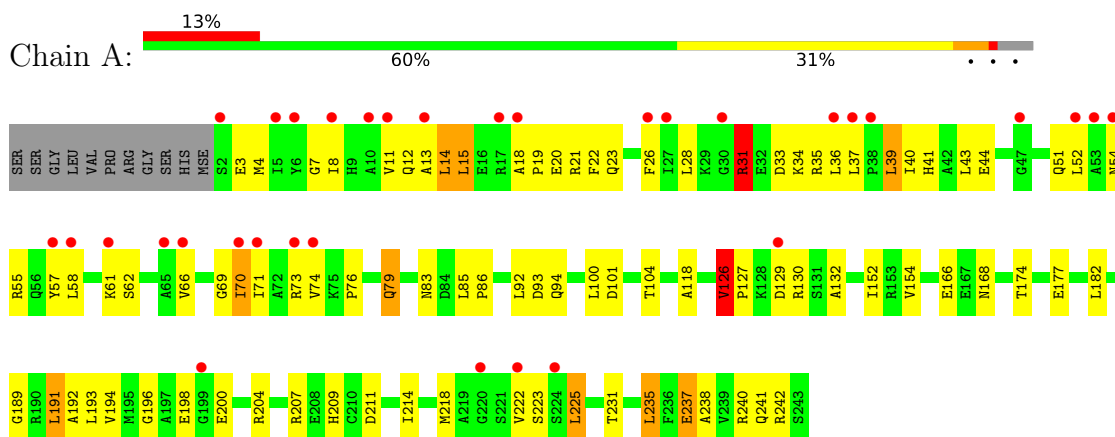
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	G	9	Total	O	0	0
			9	9		
2	H	31	Total	O	0	0
			31	31		

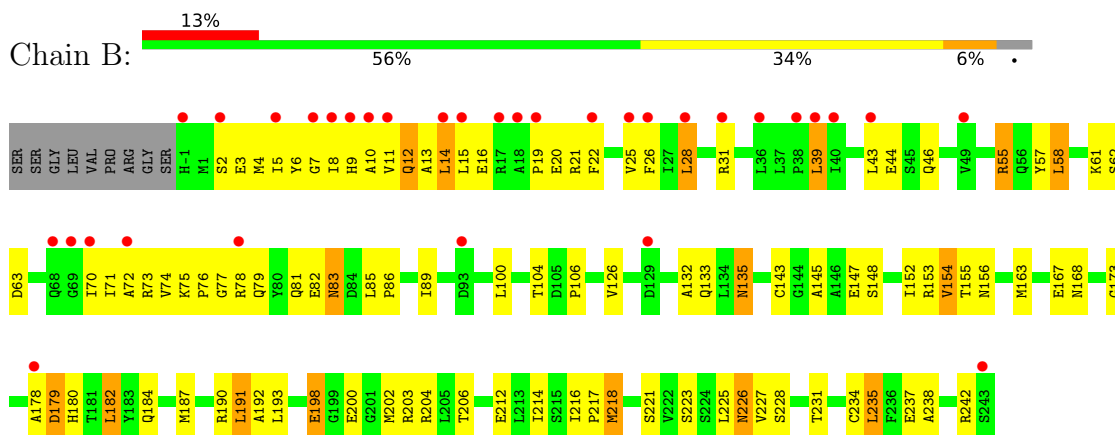
3 Residue-property plots [i](#)

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.

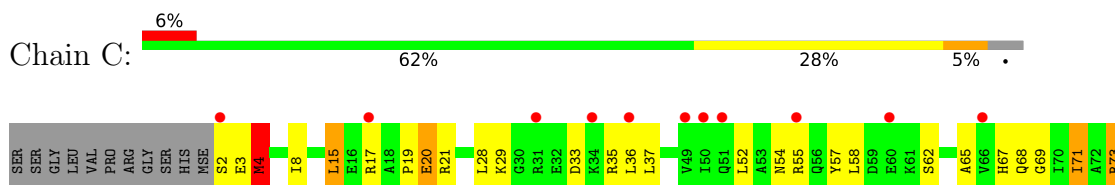
• Molecule 1: HYPOTHETICAL TRNA/RRNA METHYLTRANSFERASE YJFH

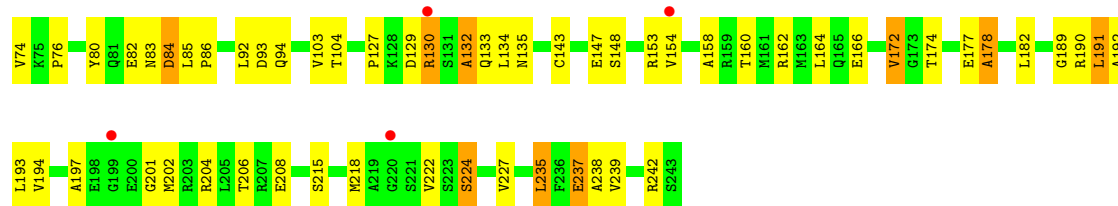


• Molecule 1: HYPOTHETICAL TRNA/RRNA METHYLTRANSFERASE YJFH

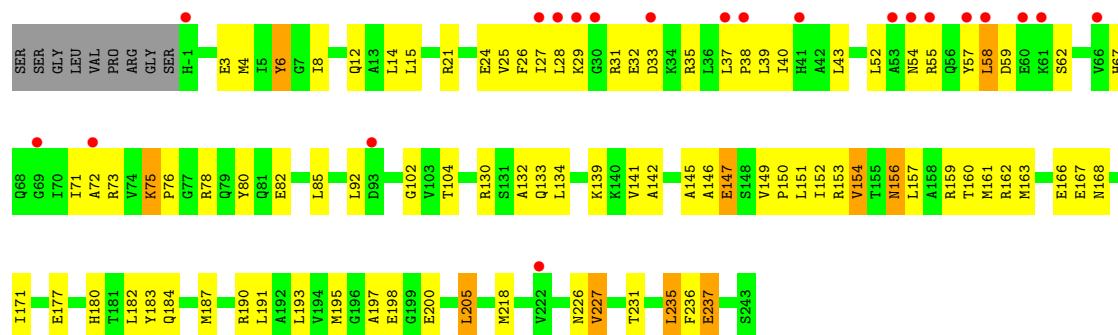


• Molecule 1: HYPOTHETICAL TRNA/RRNA METHYLTRANSFERASE YJFH

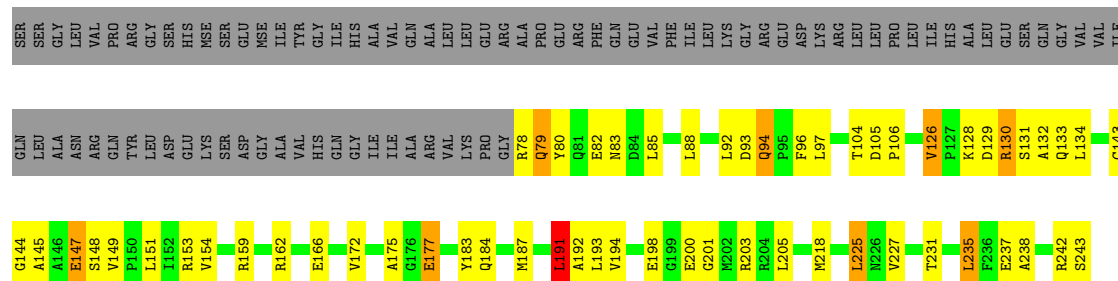




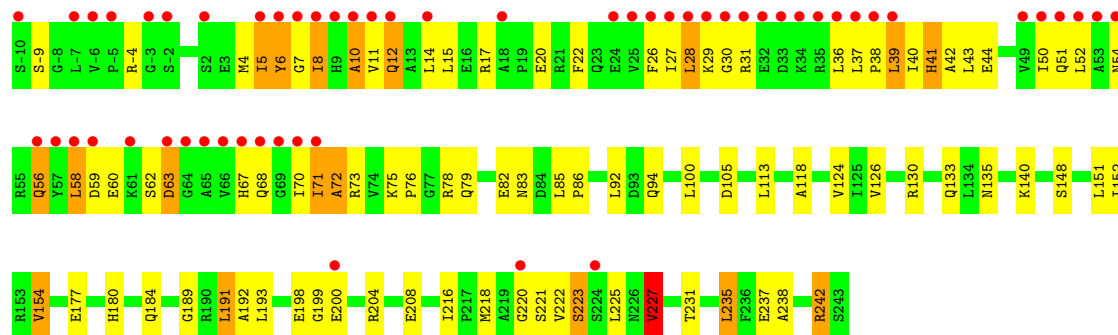
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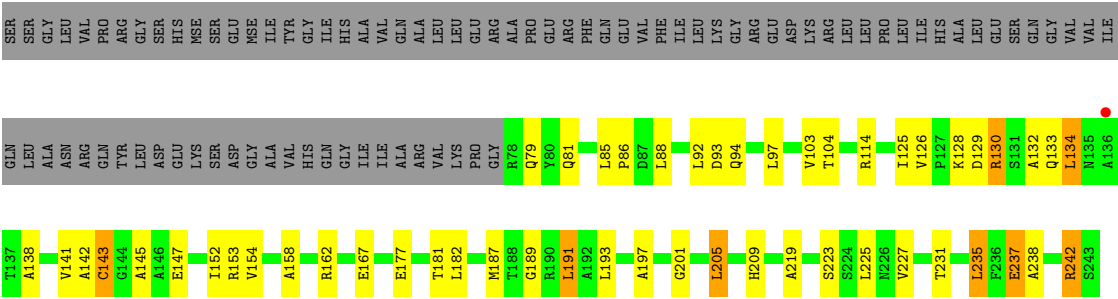


• Molecule 1: HYPOTHETICAL TRNA/RRNA METHYLTRANSFERASE YJFH



• Molecule 1: HYPOTHETICAL TRNA/RRNA METHYLTRANSFERASE YJFH

Chain G: 45% 17% . 34%



• Molecule 1: HYPOTHETICAL TRNA/RRNA METHYLTRANSFERASE YJFH

Chain H: 10% 71% 26% ..



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	87.48Å 132.29Å 90.71Å 90.00° 96.40° 90.00°	Depositor
Resolution (Å)	46.86 – 2.50 46.86 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.3 (46.86-2.50) 99.3 (46.86-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.58 (at 2.48Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.231 , 0.279 0.230 , 0.278	Depositor DCC
R_{free} test set	3587 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	43.8	Xtriage
Anisotropy	0.255	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 48.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.000 for l,-k,h	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	13953	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 55.12 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.3107e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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	0/1869	0.99	8/2514 (0.3%)
1	B	0.53	0/1887	1.00	6/2536 (0.2%)
1	C	0.57	2/1869 (0.1%)	1.03	3/2514 (0.1%)
1	D	0.55	0/1887	1.02	4/2536 (0.2%)
1	E	0.60	0/1257	1.10	9/1691 (0.5%)
1	F	0.56	0/1947	1.02	8/2617 (0.3%)
1	G	0.59	0/1257	1.08	6/1691 (0.4%)
1	H	0.54	0/1931	1.01	4/2596 (0.2%)
All	All	0.56	2/13904 (0.0%)	1.03	48/18695 (0.3%)

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
1	D	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	4	MSE	CG-SE	-5.69	1.78	1.95
1	C	4	MSE	SE-CE	-5.69	1.78	1.95

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	237	GLU	N-CA-C	-9.06	101.49	111.36
1	C	237	GLU	N-CA-C	-8.49	101.97	111.82
1	G	237	GLU	N-CA-C	-8.41	101.34	111.69
1	B	237	GLU	N-CA-C	-8.39	101.36	111.69
1	F	237	GLU	N-CA-C	-7.96	102.59	111.82
1	E	149	VAL	N-CA-C	7.72	116.53	107.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	56	GLN	N-CA-C	-7.27	103.28	111.14
1	A	191	LEU	N-CA-C	7.10	120.72	109.50
1	E	191	LEU	N-CA-C	6.95	119.99	109.23
1	D	75	LYS	N-CA-C	-6.92	100.94	109.65
1	C	227	VAL	N-CA-C	6.91	117.66	110.62
1	E	237	GLU	N-CA-C	-6.60	103.91	112.23
1	H	191	LEU	N-CA-C	6.57	119.42	109.23
1	E	238	ALA	N-CA-C	-6.46	104.16	111.14
1	D	191	LEU	N-CA-C	6.33	119.50	109.50
1	F	216	ILE	N-CA-C	-6.21	100.81	107.77
1	C	191	LEU	N-CA-C	6.16	118.78	109.23
1	F	227	VAL	CB-CA-C	-6.15	103.98	112.04
1	G	191	LEU	N-CA-C	6.06	119.37	109.24
1	F	7	GLY	N-CA-C	6.06	120.30	112.54
1	G	227	VAL	N-CA-C	6.02	116.76	110.62
1	F	191	LEU	N-CA-C	5.95	118.32	108.99
1	A	237	GLU	N-CA-C	-5.93	104.81	111.28
1	A	39	LEU	N-CA-C	-5.91	105.15	112.72
1	A	168	ASN	N-CA-C	5.73	119.57	112.24
1	F	8	ILE	N-CA-C	-5.68	103.95	111.05
1	G	181	THR	N-CA-C	5.67	117.84	110.53
1	H	237	GLU	N-CA-C	-5.59	105.34	111.82
1	B	77	GLY	N-CA-C	-5.58	105.39	112.54
1	H	126	VAL	N-CA-C	5.58	113.77	109.19
1	A	166	GLU	N-CA-C	-5.56	106.34	113.01
1	D	32	GLU	N-CA-C	-5.54	103.59	110.41
1	G	103	VAL	N-CA-C	-5.51	100.38	108.54
1	E	201	GLY	N-CA-C	-5.51	104.14	112.84
1	F	6	TYR	N-CA-C	5.49	122.49	110.80
1	E	184	GLN	N-CA-C	-5.42	105.98	112.59
1	B	191	LEU	N-CA-C	5.38	118.01	109.50
1	H	227	VAL	N-CA-C	5.38	117.77	111.05
1	A	198	GLU	N-CA-C	5.33	117.84	111.71
1	B	62	SER	N-CA-C	-5.29	106.72	114.39
1	A	70	ILE	N-CA-C	5.27	115.86	108.17
1	E	194	VAL	N-CA-C	5.25	115.42	107.75
1	B	178	ALA	N-CA-C	5.23	118.09	109.72
1	G	126	VAL	CB-CA-C	-5.21	104.23	109.33
1	B	198	GLU	N-CA-C	5.10	117.50	111.33
1	E	203	ARG	N-CA-C	-5.08	103.34	110.50
1	A	126	VAL	N-CA-C	5.02	113.31	109.19
1	E	130	ARG	NE-CZ-NH2	5.01	123.71	119.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	183	TYR	Sidechain

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	1851	0	1887	82	0
1	B	1869	0	1906	96	0
1	C	1851	0	1887	71	0
1	D	1869	0	1906	68	0
1	E	1248	0	1264	33	0
1	F	1928	0	1967	96	0
1	G	1248	0	1264	36	0
1	H	1912	0	1954	47	0
2	A	14	0	0	0	0
2	B	19	0	0	1	0
2	C	14	0	0	0	0
2	D	24	0	0	0	0
2	E	28	0	0	1	0
2	F	38	0	0	0	0
2	G	9	0	0	0	0
2	H	31	0	0	0	0
All	All	13953	0	14035	514	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 18.

All (514) 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:A:79:GLN:HE21	1:A:79:GLN:HA	1.24	1.00
1:D:85:LEU:HD11	1:D:152:ILE:HG21	1.47	0.97
1:F:26:PHE:HB2	1:F:71:ILE:HG13	1.49	0.94
1:C:8:ILE:HD11	1:C:36:LEU:HD23	1.50	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:12:GLN:HA	1:F:12:GLN:HE21	1.33	0.93
1:A:3:GLU:HG3	1:A:76:PRO:HG3	1.47	0.93
1:B:25:VAL:HG21	1:B:43:LEU:HD13	1.49	0.92
1:A:52:LEU:HD23	1:A:52:LEU:H	1.36	0.90
1:B:15:LEU:HD21	1:B:43:LEU:HD23	1.52	0.89
1:C:172:VAL:HG23	1:C:191:LEU:HD11	1.56	0.88
1:A:118:ALA:HA	1:F:223:SER:HB3	1.53	0.88
1:F:58:LEU:HD22	1:F:71:ILE:HD11	1.56	0.87
1:B:20:GLU:CD	1:B:20:GLU:H	1.82	0.86
1:A:130:ARG:HG3	1:A:130:ARG:HH11	1.41	0.84
1:A:28:LEU:HD13	1:A:69:GLY:HA2	1.57	0.84
1:G:187:MSE:HE2	1:G:191:LEU:HD21	1.57	0.84
1:B:9:HIS:CE1	1:G:223:SER:HB3	2.12	0.84
1:C:4:MSE:HE3	1:C:73:ARG:HG3	1.60	0.83
1:F:5:ILE:HG22	1:F:72:ALA:HB3	1.61	0.81
1:A:28:LEU:HD12	1:A:58:LEU:HD12	1.63	0.81
1:G:182:LEU:HG	1:G:237:GLU:HG2	1.63	0.81
1:B:28:LEU:HD22	1:B:31:ARG:HB2	1.64	0.79
1:A:79:GLN:HE21	1:A:79:GLN:CA	1.95	0.79
1:E:187:MSE:HE2	1:E:191:LEU:HD21	1.65	0.79
1:B:4:MSE:HG3	1:B:72:ALA:O	1.81	0.79
1:G:88:LEU:O	1:G:92:LEU:HD13	1.83	0.79
1:D:85:LEU:CD1	1:D:152:ILE:HG21	2.12	0.79
1:A:28:LEU:HB2	1:A:31:ARG:HD2	1.64	0.78
1:C:29:LYS:HB2	1:C:52:LEU:HB3	1.66	0.78
1:E:200:GLU:HG3	2:E:2016:HOH:O	1.82	0.77
1:F:29:LYS:O	1:F:31:ARG:HG3	1.84	0.76
1:B:85:LEU:HD11	1:B:152:ILE:HD13	1.67	0.76
1:A:129:ASP:OD1	1:D:168:ASN:HB3	1.86	0.75
1:F:82:GLU:HG3	1:F:154:VAL:HG22	1.68	0.75
1:B:58:LEU:HD22	1:B:71:ILE:CD1	2.16	0.75
1:B:85:LEU:HD13	1:B:152:ILE:HG21	1.70	0.74
1:H:21:ARG:HH12	1:H:147:GLU:HG3	1.50	0.74
1:A:3:GLU:OE1	1:A:76:PRO:HA	1.89	0.73
1:C:58:LEU:HD22	1:C:71:ILE:HG13	1.70	0.73
1:B:180:HIS:HB2	1:B:184:GLN:HB2	1.69	0.73
1:H:33:ASP:HB3	1:H:36:LEU:HB2	1.71	0.73
1:C:17:ARG:HG3	1:C:148:SER:CB	2.19	0.73
1:F:238:ALA:O	1:F:242:ARG:HG2	1.89	0.72
1:D:133:GLN:HA	1:D:153:ARG:NH2	2.04	0.72
1:A:23:GLN:HB3	1:A:73:ARG:HH21	1.55	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:218:MSE:CE	1:C:224:SER:HA	2.18	0.72
1:B:5:ILE:HD13	1:B:14:LEU:HD12	1.70	0.72
1:F:94:GLN:NE2	1:F:189:GLY:HA2	2.04	0.72
1:A:126:VAL:HG22	1:A:127:PRO:HD2	1.72	0.71
1:F:40:ILE:O	1:F:44:GLU:HG3	1.90	0.71
1:A:15:LEU:O	1:A:19:PRO:HG3	1.91	0.71
1:A:193:LEU:HD23	1:A:235:LEU:HD13	1.73	0.71
1:B:135:ASN:H	1:B:135:ASN:HD22	1.39	0.71
1:H:21:ARG:NH1	1:H:147:GLU:HG3	2.06	0.71
1:G:187:MSE:HE1	1:G:193:LEU:HD13	1.74	0.70
1:A:31:ARG:O	1:A:31:ARG:HD3	1.91	0.70
1:B:81:GLN:HE21	1:B:81:GLN:HA	1.56	0.70
1:B:238:ALA:O	1:B:242:ARG:HG3	1.92	0.69
1:F:27:ILE:HD13	1:F:40:ILE:HD13	1.75	0.69
1:C:218:MSE:HE3	1:C:224:SER:HA	1.74	0.69
1:H:85:LEU:HB3	1:H:86:PRO:HD3	1.73	0.69
1:F:222:VAL:HG23	1:F:223:SER:H	1.58	0.69
1:F:78:ARG:HG2	1:F:78:ARG:HH11	1.59	0.68
1:C:172:VAL:CG2	1:C:191:LEU:HD11	2.24	0.68
1:C:62:SER:HB3	1:C:71:ILE:HD13	1.75	0.68
1:A:118:ALA:HA	1:F:223:SER:CB	2.25	0.67
1:D:85:LEU:HD11	1:D:152:ILE:HD13	1.76	0.67
1:B:85:LEU:CD1	1:B:152:ILE:HD13	2.24	0.67
1:E:187:MSE:HE1	1:E:193:LEU:HD13	1.76	0.67
1:C:17:ARG:HG3	1:C:148:SER:HB3	1.76	0.67
1:H:182:LEU:HD23	1:H:182:LEU:H	1.60	0.67
1:F:85:LEU:HD13	1:F:152:ILE:HD13	1.75	0.66
1:B:135:ASN:HD22	1:B:135:ASN:N	1.91	0.66
1:E:198:GLU:HB2	1:E:227:VAL:HG23	1.77	0.66
1:A:40:ILE:O	1:A:44:GLU:HG3	1.95	0.65
1:B:182:LEU:HB2	1:B:214:ILE:HD12	1.77	0.65
1:F:12:GLN:HA	1:F:12:GLN:NE2	2.09	0.65
1:C:197:ALA:O	1:C:201:GLY:HA2	1.96	0.65
1:F:29:LYS:HA	1:F:52:LEU:HB3	1.78	0.65
1:B:9:HIS:NE2	1:G:223:SER:HB3	2.11	0.65
1:A:238:ALA:O	1:A:242:ARG:HG3	1.97	0.65
1:C:204:ARG:O	1:C:208:GLU:HG3	1.97	0.65
1:F:40:ILE:HG23	1:F:50:ILE:HD13	1.79	0.65
1:A:28:LEU:HD13	1:A:69:GLY:CA	2.28	0.64
1:B:14:LEU:CD1	1:B:74:VAL:HG21	2.27	0.64
1:D:28:LEU:HD23	1:D:58:LEU:HD12	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:159:ARG:HA	1:E:162:ARG:NH1	2.13	0.64
1:A:218:MSE:HE1	1:A:225:LEU:HD22	1.78	0.64
1:B:2:SER:HB3	1:B:73:ARG:HG2	1.80	0.64
1:D:133:GLN:HA	1:D:153:ARG:HH22	1.62	0.64
1:E:218:MSE:HE1	1:E:225:LEU:HD22	1.80	0.64
1:D:134:LEU:HB3	1:D:139:LYS:NZ	2.14	0.63
1:F:56:GLN:O	1:F:60:GLU:HG3	1.99	0.63
1:A:14:LEU:HB3	1:A:22:PHE:HE1	1.64	0.63
1:C:85:LEU:HB3	1:C:86:PRO:HD3	1.80	0.63
1:F:27:ILE:HG13	1:F:52:LEU:HD23	1.81	0.63
1:D:4:MSE:HE2	1:D:71:ILE:HG22	1.81	0.63
1:D:21:ARG:HD2	1:D:75:LYS:O	1.99	0.63
1:D:4:MSE:HE3	1:D:72:ALA:C	2.24	0.62
1:A:28:LEU:CB	1:A:31:ARG:HD2	2.28	0.62
1:H:88:LEU:O	1:H:92:LEU:HD22	1.99	0.62
1:D:8:ILE:HD13	1:D:35:ARG:HG2	1.81	0.62
1:H:240:ARG:HH11	1:H:240:ARG:HB3	1.64	0.62
1:C:28:LEU:HD12	1:C:28:LEU:O	1.99	0.62
1:F:82:GLU:HG3	1:F:154:VAL:CG2	2.29	0.62
1:F:58:LEU:HA	1:F:71:ILE:HD11	1.80	0.62
1:F:75:LYS:HB3	1:F:76:PRO:HD2	1.82	0.62
1:D:6:TYR:HD2	1:D:71:ILE:HG12	1.65	0.61
1:E:128:LYS:HG3	1:E:154:VAL:O	2.00	0.61
1:B:14:LEU:HD11	1:B:74:VAL:HG21	1.83	0.61
1:B:133:GLN:HA	1:B:153:ARG:NH2	2.16	0.61
1:C:193:LEU:HD23	1:C:235:LEU:HD13	1.83	0.61
1:D:85:LEU:CD1	1:D:152:ILE:HD13	2.31	0.61
1:H:88:LEU:HD21	1:H:123:ALA:HB2	1.83	0.60
1:A:14:LEU:HD21	1:A:74:VAL:HG11	1.82	0.60
1:C:2:SER:HA	1:C:76:PRO:HD3	1.82	0.60
1:B:218:MSE:HG2	1:B:223:SER:HB2	1.82	0.60
1:E:193:LEU:HD23	1:E:235:LEU:HD13	1.82	0.60
1:A:126:VAL:HG22	1:A:127:PRO:CD	2.31	0.60
1:H:27:ILE:HG13	1:H:52:LEU:HG	1.83	0.60
1:A:33:ASP:C	1:A:35:ARG:H	2.09	0.60
1:A:130:ARG:HG3	1:A:130:ARG:NH1	2.12	0.60
1:B:15:LEU:HD21	1:B:43:LEU:CD2	2.29	0.59
1:B:218:MSE:HG2	1:B:223:SER:CB	2.31	0.59
2:B:2004:HOH:O	1:G:219:ALA:HB2	2.01	0.59
1:A:28:LEU:CD1	1:A:69:GLY:HA2	2.30	0.59
1:F:59:ASP:HA	1:F:62:SER:OG	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:25:VAL:HG21	1:D:43:LEU:HD13	1.84	0.59
1:F:36:LEU:O	1:F:40:ILE:HG13	2.03	0.59
1:A:11:VAL:HG11	1:A:70:ILE:HD11	1.85	0.59
1:B:82:GLU:CD	1:B:154:VAL:HG22	2.28	0.59
1:C:33:ASP:O	1:C:37:LEU:HD13	2.03	0.59
1:F:29:LYS:HA	1:F:52:LEU:HD13	1.85	0.59
1:B:12:GLN:C	1:B:14:LEU:H	2.11	0.58
1:B:28:LEU:H	1:B:28:LEU:HD12	1.68	0.58
1:B:135:ASN:H	1:B:135:ASN:ND2	2.01	0.58
1:G:142:ALA:HB1	1:G:145:ALA:HB3	1.86	0.58
1:B:28:LEU:HD12	1:B:28:LEU:N	2.17	0.58
1:H:193:LEU:HD23	1:H:235:LEU:HD13	1.84	0.58
1:D:157:LEU:O	1:D:161:MSE:HG3	2.04	0.58
1:F:58:LEU:HD22	1:F:71:ILE:CD1	2.31	0.58
1:D:62:SER:HB2	1:D:71:ILE:HD11	1.85	0.58
1:F:26:PHE:HB2	1:F:71:ILE:CG1	2.29	0.57
1:F:130:ARG:NH2	1:F:200:GLU:HG3	2.19	0.57
1:A:85:LEU:N	1:A:86:PRO:HD2	2.19	0.57
1:A:4:MSE:CE	1:A:73:ARG:HG3	2.35	0.57
1:D:58:LEU:HD23	1:D:71:ILE:HD12	1.86	0.57
1:D:180:HIS:HB2	1:D:184:GLN:OE1	2.03	0.57
1:B:11:VAL:HG21	1:B:70:ILE:HD13	1.87	0.57
1:B:82:GLU:OE2	1:B:154:VAL:HG22	2.05	0.57
1:C:21:ARG:O	1:C:74:VAL:HG13	2.04	0.57
1:G:85:LEU:N	1:G:86:PRO:HD2	2.20	0.57
1:H:62:SER:O	1:H:63:ASP:HB3	2.05	0.57
1:H:203:ARG:HG2	1:H:203:ARG:HH11	1.69	0.57
1:G:129:ASP:O	1:G:130:ARG:HB2	2.05	0.56
1:G:205:LEU:HD23	1:G:209:HIS:CE1	2.40	0.56
1:G:177:GLU:CD	1:G:177:GLU:H	2.13	0.56
1:H:28:LEU:HD22	1:H:31:ARG:HB2	1.86	0.56
1:A:37:LEU:HD12	1:A:40:ILE:HD12	1.88	0.56
1:B:78:ARG:HH11	1:B:78:ARG:HG2	1.70	0.56
1:C:15:LEU:O	1:C:19:PRO:HG3	2.06	0.56
1:F:12:GLN:HE21	1:F:12:GLN:CA	2.09	0.56
1:E:133:GLN:HA	1:E:153:ARG:NH2	2.19	0.56
1:D:130:ARG:HG3	1:D:130:ARG:HH11	1.71	0.56
1:E:80:TYR:CE2	1:E:88:LEU:HD13	2.41	0.56
1:B:231:THR:HG22	1:B:235:LEU:HD22	1.88	0.56
1:G:128:LYS:HG2	1:G:153:ARG:HB3	1.89	0.56
1:H:133:GLN:HA	1:H:153:ARG:NH2	2.19	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:231:THR:O	1:B:235:LEU:HB2	2.05	0.55
1:F:28:LEU:HD22	1:F:31:ARG:HB2	1.88	0.55
1:G:205:LEU:HD23	1:G:209:HIS:HE1	1.71	0.55
1:A:85:LEU:HD13	1:A:152:ILE:HD13	1.87	0.55
1:A:193:LEU:CD2	1:A:235:LEU:HD13	2.35	0.55
1:H:231:THR:HG22	1:H:235:LEU:HD22	1.87	0.55
1:A:52:LEU:H	1:A:52:LEU:CD2	2.15	0.55
1:F:-4:ARG:HH11	1:F:-4:ARG:HG2	1.72	0.55
1:A:4:MSE:HE1	1:A:73:ARG:HG3	1.88	0.55
1:A:100:LEU:O	1:A:127:PRO:HD3	2.06	0.55
1:H:15:LEU:HA	1:H:22:PHE:HE2	1.71	0.55
1:A:204:ARG:HA	1:A:207:ARG:NH1	2.21	0.55
1:B:81:GLN:HA	1:B:81:GLN:NE2	2.21	0.55
1:B:191:LEU:HD12	1:B:192:ALA:H	1.71	0.55
1:F:100:LEU:HB2	1:F:126:VAL:HG12	1.88	0.55
1:A:13:ALA:HA	1:F:222:VAL:CG1	2.37	0.55
1:A:85:LEU:HD13	1:A:152:ILE:CD1	2.37	0.55
1:H:37:LEU:HB2	1:H:38:PRO:HD3	1.89	0.55
1:D:78:ARG:HH11	1:D:78:ARG:HG2	1.72	0.55
1:D:236:PHE:HB3	1:E:183:TYR:OH	2.07	0.55
1:D:104:THR:O	1:D:132:ALA:HB2	2.08	0.54
1:F:39:LEU:C	1:F:41:HIS:H	2.15	0.54
1:A:7:GLY:O	1:A:11:VAL:HG23	2.07	0.54
1:C:3:GLU:OE2	1:C:76:PRO:HA	2.07	0.54
1:H:82:GLU:HG3	1:H:154:VAL:HG22	1.89	0.54
1:C:194:VAL:HG12	1:C:202:MSE:HE2	1.90	0.54
1:A:37:LEU:C	1:A:39:LEU:H	2.16	0.54
1:B:4:MSE:HE2	1:B:73:ARG:HG3	1.90	0.54
1:H:104:THR:C	1:H:132:ALA:HB2	2.33	0.54
1:C:21:ARG:HB3	1:C:74:VAL:CG1	2.38	0.54
1:C:127:PRO:HG3	1:C:130:ARG:NH1	2.23	0.54
1:C:147:GLU:HA	1:C:147:GLU:OE1	2.07	0.53
1:F:218:MSE:HB3	1:F:223:SER:HB2	1.90	0.53
1:B:180:HIS:HE2	1:B:212:GLU:HG3	1.74	0.53
1:C:133:GLN:HA	1:C:153:ARG:NH2	2.24	0.53
1:F:15:LEU:HD12	1:F:43:LEU:HG	1.90	0.53
1:H:128:LYS:HG2	1:H:153:ARG:HB3	1.89	0.53
1:A:21:ARG:HB3	1:A:74:VAL:CG1	2.38	0.53
1:C:84:ASP:N	1:C:84:ASP:OD1	2.41	0.53
1:F:17:ARG:HG2	1:F:17:ARG:HH11	1.73	0.53
1:F:58:LEU:CD2	1:F:71:ILE:HD11	2.35	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:28:LEU:CD1	1:A:58:LEU:HD12	2.36	0.53
1:D:28:LEU:HD23	1:D:58:LEU:CD1	2.39	0.53
1:A:28:LEU:HD21	1:A:55:ARG:HH21	1.74	0.53
1:E:126:VAL:HG13	1:E:131:SER:OG	2.09	0.52
1:F:44:GLU:HG2	1:F:50:ILE:HD12	1.91	0.52
1:A:58:LEU:HD22	1:A:71:ILE:HD12	1.89	0.52
1:F:78:ARG:HG2	1:F:78:ARG:NH1	2.24	0.52
1:G:197:ALA:O	1:G:201:GLY:HA2	2.10	0.52
1:B:58:LEU:HD22	1:B:71:ILE:HD11	1.90	0.52
1:B:83:ASN:O	1:B:86:PRO:HD2	2.10	0.52
1:C:8:ILE:HD11	1:C:36:LEU:CD2	2.32	0.52
1:D:141:VAL:HG12	1:D:141:VAL:O	2.08	0.52
1:F:29:LYS:CA	1:F:52:LEU:HB3	2.39	0.52
1:G:187:MSE:O	1:G:242:ARG:HD2	2.09	0.52
1:A:240:ARG:NH1	1:A:241:GLN:HG3	2.24	0.52
1:C:54:ASN:HD21	1:C:57:TYR:HB2	1.74	0.52
1:F:113:LEU:HD23	1:F:124:VAL:HG21	1.91	0.52
1:G:242:ARG:HG2	1:G:242:ARG:NH1	2.24	0.52
1:H:240:ARG:HH12	1:H:241:GLN:HG3	1.74	0.52
1:G:231:THR:HG22	1:G:235:LEU:HD22	1.92	0.52
1:C:3:GLU:HG3	1:C:76:PRO:HG3	1.92	0.52
1:D:134:LEU:HB3	1:D:139:LYS:HZ2	1.75	0.52
1:E:94:GLN:HE21	1:E:94:GLN:CA	2.23	0.52
1:B:20:GLU:CD	1:B:20:GLU:N	2.60	0.51
1:B:21:ARG:HA	1:B:75:LYS:HG2	1.92	0.51
1:G:231:THR:O	1:G:235:LEU:HB2	2.10	0.51
1:C:52:LEU:HD12	1:C:52:LEU:N	2.25	0.51
1:C:133:GLN:HA	1:C:153:ARG:HH22	1.75	0.51
1:F:6:TYR:HE2	1:F:62:SER:HA	1.75	0.51
1:F:58:LEU:HA	1:F:71:ILE:CD1	2.41	0.51
1:B:12:GLN:O	1:B:16:GLU:HG2	2.11	0.51
1:H:4:MSE:HE3	1:H:71:ILE:HG22	1.91	0.51
1:C:33:ASP:C	1:C:37:LEU:HD13	2.36	0.51
1:H:52:LEU:HD12	1:H:52:LEU:N	2.26	0.51
1:A:118:ALA:CA	1:F:223:SER:HB3	2.35	0.51
1:A:11:VAL:HG11	1:A:70:ILE:CD1	2.41	0.51
1:E:231:THR:HG22	1:E:235:LEU:HD22	1.93	0.51
1:C:17:ARG:HG3	1:C:148:SER:HB2	1.90	0.51
1:C:235:LEU:O	1:C:238:ALA:HB3	2.11	0.51
1:F:83:ASN:O	1:F:86:PRO:HD2	2.11	0.51
1:A:40:ILE:HA	1:A:43:LEU:HD12	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:174:THR:HA	1:A:214:ILE:O	2.11	0.51
1:F:37:LEU:N	1:F:38:PRO:HD2	2.26	0.50
1:D:27:ILE:HG21	1:D:40:ILE:HD11	1.94	0.50
1:B:21:ARG:HD2	1:B:75:LYS:HG3	1.93	0.50
1:B:106:PRO:HB3	1:B:135:ASN:HD21	1.76	0.50
1:D:182:LEU:HG	1:D:237:GLU:HG2	1.92	0.50
1:F:82:GLU:CG	1:F:154:VAL:HG22	2.40	0.50
1:A:3:GLU:CG	1:A:76:PRO:HG3	2.32	0.50
1:D:193:LEU:HD23	1:D:235:LEU:HD13	1.91	0.50
1:F:40:ILE:O	1:F:40:ILE:HG22	2.11	0.50
1:F:26:PHE:HD2	1:F:71:ILE:HB	1.77	0.50
1:F:28:LEU:HD13	1:F:31:ARG:HB2	1.93	0.50
1:A:21:ARG:O	1:A:74:VAL:HG13	2.12	0.50
1:F:12:GLN:HE22	1:F:15:LEU:HD13	1.76	0.50
1:B:216:ILE:HD11	1:B:234:CYS:SG	2.51	0.50
1:D:146:ALA:HB3	1:D:147:GLU:OE2	2.11	0.50
1:H:96:PHE:C	1:H:97:LEU:HD23	2.37	0.50
1:H:7:GLY:HA2	1:H:67:HIS:O	2.11	0.50
1:B:145:ALA:HA	1:B:148:SER:OG	2.12	0.49
1:B:21:ARG:HB3	1:B:74:VAL:HG12	1.93	0.49
1:E:94:GLN:HE21	1:E:94:GLN:HA	1.77	0.49
1:F:15:LEU:HD23	1:F:15:LEU:C	2.37	0.49
1:G:104:THR:O	1:G:132:ALA:HB2	2.13	0.49
1:C:104:THR:C	1:C:132:ALA:HB2	2.38	0.49
1:C:178:ALA:HB3	1:C:215:SER:HB3	1.95	0.49
1:B:193:LEU:HD23	1:B:235:LEU:HD13	1.94	0.49
1:D:130:ARG:HH11	1:D:130:ARG:CG	2.25	0.49
1:D:102:GLY:O	1:D:197:ALA:HB2	2.13	0.49
1:E:191:LEU:HB3	1:E:242:ARG:NH2	2.28	0.49
1:C:160:THR:O	1:C:164:LEU:HG	2.12	0.49
1:C:218:MSE:HE1	1:C:224:SER:HA	1.93	0.49
1:F:67:HIS:O	1:F:68:GLN:HB2	2.11	0.49
1:G:104:THR:C	1:G:132:ALA:HB2	2.38	0.49
1:H:218:MSE:HE2	1:H:218:MSE:HA	1.94	0.49
1:B:57:TYR:CZ	1:B:61:LYS:HD2	2.48	0.49
1:C:104:THR:O	1:C:132:ALA:HB2	2.13	0.49
1:F:8:ILE:O	1:F:12:GLN:HB3	2.13	0.49
1:F:30:GLY:H	1:F:52:LEU:HD13	1.77	0.49
1:A:21:ARG:HB3	1:A:74:VAL:HG13	1.94	0.48
1:B:26:PHE:HB2	1:B:71:ILE:CG1	2.42	0.48
1:E:145:ALA:HA	1:E:148:SER:HB3	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:211:ASP:OD1	1:B:203:ARG:HB2	2.14	0.48
1:E:78:ARG:HG2	1:E:79:GLN:N	2.29	0.48
1:G:238:ALA:O	1:G:242:ARG:HB2	2.13	0.48
1:B:6:TYR:O	1:B:10:ALA:HB3	2.14	0.48
1:B:21:ARG:HB3	1:B:74:VAL:CG1	2.43	0.48
1:E:129:ASP:C	1:E:131:SER:H	2.21	0.48
1:F:70:ILE:HG13	1:F:70:ILE:O	2.13	0.48
1:B:25:VAL:HG21	1:B:43:LEU:CD1	2.33	0.48
1:E:104:THR:C	1:E:132:ALA:HB2	2.39	0.48
1:F:204:ARG:O	1:F:208:GLU:HG3	2.13	0.48
1:G:85:LEU:HD13	1:G:152:ILE:HD13	1.95	0.48
1:C:28:LEU:HD23	1:C:69:GLY:HA2	1.96	0.47
1:H:67:HIS:O	1:H:68:GLN:HB2	2.14	0.47
1:A:18:ALA:HB3	1:A:21:ARG:HD3	1.95	0.47
1:H:15:LEU:HD13	1:H:43:LEU:HD23	1.95	0.47
1:A:79:GLN:HA	1:A:79:GLN:NE2	2.08	0.47
1:G:242:ARG:HG2	1:G:242:ARG:HH11	1.80	0.47
1:A:8:ILE:O	1:A:12:GLN:HG3	2.15	0.47
1:C:20:GLU:H	1:C:20:GLU:HG3	1.33	0.47
1:D:37:LEU:N	1:D:38:PRO:HD2	2.29	0.47
1:D:134:LEU:HG	1:D:151:LEU:CD2	2.44	0.47
1:F:27:ILE:CD1	1:F:40:ILE:HD13	2.43	0.47
1:G:193:LEU:HD23	1:G:235:LEU:HD13	1.95	0.47
1:C:62:SER:HB2	1:C:65:ALA:HB3	1.96	0.47
1:A:31:ARG:NH1	1:A:36:LEU:HD11	2.29	0.47
1:B:198:GLU:HB2	1:B:227:VAL:HG23	1.97	0.47
1:B:198:GLU:CD	1:G:114:ARG:HH22	2.21	0.47
1:C:182:LEU:HG	1:C:237:GLU:HG2	1.97	0.47
1:D:29:LYS:C	1:D:31:ARG:H	2.22	0.47
1:D:12:GLN:HB2	1:D:39:LEU:HD21	1.95	0.47
1:F:15:LEU:HD21	1:F:42:ALA:HB1	1.96	0.47
1:F:231:THR:O	1:F:235:LEU:HB2	2.14	0.47
1:D:59:ASP:CG	1:D:67:HIS:HE2	2.23	0.47
1:A:13:ALA:HA	1:F:222:VAL:HG11	1.96	0.47
1:B:4:MSE:HE2	1:B:73:ARG:CG	2.45	0.47
1:B:78:ARG:HG2	1:B:78:ARG:NH1	2.30	0.47
1:C:202:MSE:HG2	1:C:206:THR:HB	1.97	0.47
1:C:158:ALA:O	1:C:162:ARG:HG3	2.16	0.46
1:G:158:ALA:HB1	1:G:162:ARG:HH21	1.80	0.46
1:B:19:PRO:HA	1:B:22:PHE:CD1	2.50	0.46
1:B:100:LEU:HB2	1:B:126:VAL:HG12	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:14:LEU:HD21	1:D:21:ARG:HB2	1.97	0.46
1:F:39:LEU:C	1:F:41:HIS:N	2.73	0.46
1:F:8:ILE:C	1:F:10:ALA:H	2.22	0.46
1:A:101:ASP:O	1:A:196:GLY:HA2	2.16	0.46
1:F:26:PHE:HA	1:F:51:GLN:O	2.16	0.46
1:A:191:LEU:HG	1:A:192:ALA:N	2.29	0.46
1:B:198:GLU:HB2	1:B:227:VAL:CG2	2.44	0.46
1:B:11:VAL:HG12	1:B:39:LEU:HD11	1.98	0.46
1:B:173:GLY:HA3	1:B:202:MSE:HE1	1.98	0.46
1:C:8:ILE:CD1	1:C:36:LEU:HD23	2.33	0.46
1:F:28:LEU:HB2	1:F:31:ARG:HD2	1.97	0.46
1:F:71:ILE:HG22	1:F:72:ALA:H	1.81	0.46
1:H:22:PHE:HA	1:H:74:VAL:HG12	1.98	0.46
1:B:85:LEU:CD1	1:B:152:ILE:HG21	2.41	0.46
1:B:226:ASN:OD1	1:B:228:SER:HB2	2.16	0.46
1:D:78:ARG:HH11	1:D:78:ARG:CG	2.28	0.46
1:H:15:LEU:HD12	1:H:19:PRO:HB3	1.98	0.46
1:D:6:TYR:CD2	1:D:71:ILE:HG12	2.48	0.45
1:E:134:LEU:HD12	1:E:151:LEU:HD23	1.98	0.45
1:D:27:ILE:HG13	1:D:52:LEU:HD23	1.98	0.45
1:F:17:ARG:HG2	1:F:17:ARG:NH1	2.31	0.45
1:F:177:GLU:H	1:F:177:GLU:HG3	1.52	0.45
1:A:14:LEU:HB3	1:A:22:PHE:CE1	2.49	0.45
1:B:19:PRO:HA	1:B:22:PHE:HD1	1.81	0.45
1:C:193:LEU:CD2	1:C:235:LEU:HD13	2.46	0.45
1:A:13:ALA:HA	1:F:222:VAL:HG12	1.99	0.45
1:A:26:PHE:CD1	1:A:51:GLN:HB3	2.52	0.45
1:A:104:THR:C	1:A:132:ALA:HB2	2.42	0.45
1:C:33:ASP:OD1	1:C:35:ARG:HB3	2.17	0.45
1:D:54:ASN:O	1:D:57:TYR:HB3	2.16	0.45
1:B:2:SER:CB	1:B:73:ARG:HG2	2.45	0.45
1:F:27:ILE:O	1:F:52:LEU:HA	2.17	0.45
1:F:28:LEU:HD22	1:F:31:ARG:CB	2.46	0.45
1:A:57:TYR:HD1	1:A:57:TYR:O	2.00	0.45
1:D:193:LEU:CD2	1:D:235:LEU:HD13	2.45	0.45
1:E:191:LEU:HG	1:E:192:ALA:N	2.32	0.45
1:H:14:LEU:CD2	1:H:74:VAL:HG11	2.47	0.45
1:H:28:LEU:N	1:H:28:LEU:HD12	2.32	0.45
1:B:3:GLU:CD	1:B:76:PRO:HA	2.42	0.44
1:D:195:MSE:HE2	1:D:231:THR:HA	1.99	0.44
1:E:177:GLU:H	1:E:177:GLU:CD	2.26	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:94:GLN:NE2	1:A:189:GLY:HA2	2.32	0.44
1:A:33:ASP:C	1:A:35:ARG:N	2.72	0.44
1:B:5:ILE:CD1	1:B:14:LEU:HD12	2.44	0.44
1:F:14:LEU:HA	1:F:17:ARG:HB2	1.98	0.44
1:B:191:LEU:HD12	1:B:192:ALA:N	2.32	0.44
1:F:28:LEU:HD13	1:F:31:ARG:O	2.16	0.44
1:H:82:GLU:HG3	1:H:154:VAL:CG2	2.46	0.44
1:A:37:LEU:C	1:A:39:LEU:N	2.75	0.44
1:H:179:ASP:OD1	1:H:179:ASP:C	2.60	0.44
1:A:223:SER:HB3	1:F:118:ALA:HA	2.00	0.44
1:B:218:MSE:HG2	1:B:223:SER:HB3	1.98	0.44
1:D:161:MSE:CE	1:D:171:ILE:HG21	2.48	0.44
1:D:168:ASN:O	1:D:190:ARG:HD3	2.18	0.44
1:F:105:ASP:OD1	1:F:105:ASP:C	2.60	0.44
1:D:104:THR:C	1:D:132:ALA:HB2	2.42	0.44
1:C:222:VAL:HG22	1:C:222:VAL:O	2.18	0.43
1:D:205:LEU:HD23	1:D:205:LEU:HA	1.79	0.43
1:F:200:GLU:H	1:F:200:GLU:HG2	1.58	0.43
1:F:220:GLY:C	1:F:222:VAL:H	2.25	0.43
1:B:12:GLN:C	1:B:14:LEU:N	2.75	0.43
1:D:218:MSE:HA	1:D:218:MSE:HE2	2.01	0.43
1:F:-4:ARG:HG2	1:F:-4:ARG:NH1	2.33	0.43
1:F:8:ILE:HG12	1:F:70:ILE:HD13	1.99	0.43
1:A:39:LEU:HD22	1:A:43:LEU:HD11	2.00	0.43
1:G:94:GLN:NE2	1:G:189:GLY:HA2	2.34	0.43
1:G:147:GLU:OE1	1:G:147:GLU:N	2.52	0.43
1:B:7:GLY:O	1:B:11:VAL:HG23	2.18	0.43
1:C:33:ASP:C	1:C:35:ARG:N	2.76	0.43
1:D:102:GLY:HA2	1:D:130:ARG:HG2	2.00	0.43
1:D:167:GLU:O	1:D:190:ARG:NH1	2.51	0.43
1:E:82:GLU:O	1:E:85:LEU:HB2	2.19	0.43
1:F:198:GLU:HG3	1:F:227:VAL:HG22	2.00	0.43
1:G:141:VAL:C	1:G:143:CYS:N	2.77	0.43
1:A:209:HIS:HA	1:B:204:ARG:HD2	2.01	0.43
1:E:144:GLY:O	1:E:147:GLU:HG2	2.18	0.43
1:F:191:LEU:HG	1:F:192:ALA:N	2.33	0.43
1:C:135:ASN:OD1	1:C:135:ASN:C	2.62	0.43
1:D:14:LEU:HD23	1:D:14:LEU:O	2.19	0.43
1:D:78:ARG:HD3	1:D:80:TYR:CZ	2.53	0.43
1:F:180:HIS:HB2	1:F:184:GLN:OE1	2.19	0.43
1:B:44:GLU:C	1:B:46:GLN:H	2.25	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:204:ARG:HD3	1:C:208:GLU:OE2	2.19	0.43
1:D:28:LEU:HD12	1:D:31:ARG:HD3	2.01	0.43
1:H:182:LEU:HD23	1:H:182:LEU:N	2.32	0.43
1:H:182:LEU:HB3	1:H:214:ILE:HD12	2.00	0.43
1:A:37:LEU:HA	1:A:40:ILE:HD12	2.00	0.43
1:C:83:ASN:O	1:C:86:PRO:HD2	2.19	0.43
1:G:235:LEU:HD12	1:G:235:LEU:HA	1.86	0.43
1:H:51:GLN:C	1:H:52:LEU:HD12	2.44	0.43
1:D:231:THR:HG22	1:D:235:LEU:HD22	2.01	0.43
1:F:6:TYR:CE2	1:F:62:SER:HA	2.54	0.43
1:F:140:LYS:HE3	1:F:140:LYS:HB2	1.85	0.43
1:H:58:LEU:HB3	1:H:67:HIS:NE2	2.34	0.43
1:H:193:LEU:CD2	1:H:235:LEU:HD13	2.47	0.43
1:C:174:THR:H	1:C:202:MSE:HE1	1.84	0.42
1:D:4:MSE:HE3	1:D:73:ARG:N	2.33	0.42
1:B:21:ARG:HD2	1:B:75:LYS:CG	2.48	0.42
1:B:26:PHE:CD1	1:B:26:PHE:N	2.86	0.42
1:D:149:VAL:HA	1:D:150:PRO:HD3	1.87	0.42
1:E:93:ASP:OD1	1:E:93:ASP:N	2.52	0.42
1:A:23:GLN:HB3	1:A:73:ARG:NH2	2.30	0.42
1:B:85:LEU:O	1:B:89:ILE:HG13	2.18	0.42
1:B:133:GLN:HA	1:B:153:ARG:HH22	1.82	0.42
1:D:130:ARG:CG	1:D:130:ARG:NH1	2.81	0.42
1:E:231:THR:O	1:E:235:LEU:HB2	2.19	0.42
1:A:200:GLU:H	1:A:200:GLU:HG2	1.53	0.42
1:B:104:THR:C	1:B:132:ALA:HB2	2.45	0.42
1:B:167:GLU:O	1:B:190:ARG:NH1	2.52	0.42
1:C:80:TYR:HA	1:C:84:ASP:OD2	2.19	0.42
1:D:200:GLU:H	1:D:200:GLU:HG2	1.64	0.42
1:F:56:GLN:HG2	1:F:60:GLU:CD	2.44	0.42
1:C:177:GLU:O	1:C:178:ALA:O	2.37	0.42
1:C:191:LEU:HG	1:C:192:ALA:N	2.34	0.42
1:D:198:GLU:HB2	1:D:227:VAL:HG22	2.02	0.42
1:G:134:LEU:HD23	1:G:138:ALA:HB3	2.02	0.42
1:A:182:LEU:HG	1:A:237:GLU:HG2	2.00	0.42
1:A:231:THR:HG22	1:A:235:LEU:HD22	2.01	0.42
1:B:85:LEU:HB3	1:B:163:MSE:CE	2.49	0.42
1:E:175:ALA:HB1	1:E:177:GLU:OE2	2.19	0.42
1:F:62:SER:O	1:F:63:ASP:C	2.62	0.42
1:H:106:PRO:HG3	1:H:135:ASN:HD21	1.83	0.42
1:A:8:ILE:HD11	1:A:36:LEU:CD2	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:217:PRO:O	1:B:218:MSE:HE3	2.19	0.42
1:E:104:THR:O	1:E:132:ALA:HB2	2.20	0.42
1:F:220:GLY:C	1:F:222:VAL:N	2.78	0.42
1:B:85:LEU:N	1:B:86:PRO:HD2	2.34	0.42
1:C:3:GLU:CD	1:C:76:PRO:HA	2.44	0.42
1:D:160:THR:O	1:D:163:MSE:HB3	2.20	0.42
1:F:71:ILE:O	1:F:72:ALA:HB2	2.20	0.42
1:B:31:ARG:NH1	1:B:55:ARG:HE	2.18	0.42
1:B:58:LEU:HD22	1:B:71:ILE:CG1	2.50	0.42
1:B:202:MSE:SE	1:B:206:THR:HG22	2.70	0.42
1:D:82:GLU:CD	1:D:154:VAL:HG22	2.44	0.42
1:F:54:ASN:O	1:F:58:LEU:HG	2.19	0.42
1:C:67:HIS:O	1:C:68:GLN:HB2	2.20	0.42
1:E:172:VAL:CG2	1:E:191:LEU:HD11	2.50	0.41
1:H:39:LEU:C	1:H:39:LEU:HD23	2.45	0.41
1:F:133:GLN:O	1:F:135:ASN:N	2.49	0.41
1:H:62:SER:C	1:H:64:GLY:H	2.28	0.41
1:B:58:LEU:HD22	1:B:71:ILE:HG12	2.02	0.41
1:C:238:ALA:O	1:C:242:ARG:HG3	2.20	0.41
1:A:4:MSE:HE2	1:A:73:ARG:HG3	2.03	0.41
1:C:21:ARG:C	1:C:74:VAL:HG13	2.44	0.41
1:H:15:LEU:CA	1:H:22:PHE:HE2	2.32	0.41
1:B:82:GLU:CG	1:B:154:VAL:HG22	2.50	0.41
1:C:36:LEU:O	1:C:37:LEU:C	2.64	0.41
1:D:24:GLU:HG2	1:D:26:PHE:CE1	2.56	0.41
1:E:105:ASP:HA	1:E:106:PRO:HD3	1.89	0.41
1:F:11:VAL:HB	1:F:70:ILE:HD12	2.03	0.41
1:F:14:LEU:HD23	1:F:22:PHE:CE2	2.56	0.41
1:A:41:HIS:HA	1:A:44:GLU:OE1	2.21	0.41
1:C:33:ASP:O	1:C:36:LEU:N	2.51	0.41
1:C:62:SER:HB3	1:C:71:ILE:CD1	2.48	0.41
1:D:6:TYR:HD2	1:D:71:ILE:CG1	2.32	0.41
1:E:172:VAL:HG23	1:E:191:LEU:CD1	2.50	0.41
1:G:85:LEU:HD21	1:G:125:ILE:CD1	2.50	0.41
1:G:94:GLN:HE21	1:G:189:GLY:HA2	1.85	0.41
1:B:163:MSE:SE	1:B:163:MSE:C	3.14	0.41
1:C:189:GLY:O	1:C:242:ARG:NH2	2.54	0.41
1:H:205:LEU:HD23	1:H:205:LEU:HA	1.89	0.41
1:C:82:GLU:HG3	1:C:154:VAL:HG22	2.02	0.41
1:C:94:GLN:NE2	1:C:189:GLY:HA2	2.36	0.41
1:G:141:VAL:C	1:G:143:CYS:H	2.29	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:175:ALA:HB1	1:H:177:GLU:HG2	2.03	0.41
1:C:235:LEU:O	1:C:239:VAL:HG23	2.20	0.41
1:D:3:GLU:HG3	1:D:76:PRO:HB3	2.03	0.41
1:D:156:ASN:HB3	1:D:159:ARG:CB	2.51	0.41
1:D:142:ALA:HB1	1:D:145:ALA:HB3	2.03	0.40
1:H:167:GLU:O	1:H:190:ARG:NH1	2.54	0.40
1:C:54:ASN:ND2	1:C:57:TYR:HB2	2.35	0.40
1:D:141:VAL:O	1:D:141:VAL:CG1	2.69	0.40
1:E:96:PHE:HB2	1:E:242:ARG:NH1	2.36	0.40
1:C:28:LEU:HD22	1:C:55:ARG:HH11	1.86	0.40
1:F:85:LEU:HB3	1:F:86:PRO:HD3	2.04	0.40
1:B:79:GLN:NE2	1:B:79:GLN:HA	2.36	0.40
1:B:85:LEU:HD23	1:B:163:MSE:CE	2.52	0.40
1:B:155:THR:HG22	1:B:156:ASN:N	2.37	0.40
1:F:27:ILE:HG13	1:F:52:LEU:CD2	2.51	0.40
1:H:20:GLU:H	1:H:20:GLU:CD	2.28	0.40
1:A:209:HIS:HA	1:B:204:ARG:CD	2.51	0.40
1:G:85:LEU:CD1	1:G:152:ILE:HD13	2.50	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	240/253 (95%)	218 (91%)	17 (7%)	5 (2%)	5 9
1	B	242/253 (96%)	215 (89%)	24 (10%)	3 (1%)	10 20
1	C	240/253 (95%)	221 (92%)	15 (6%)	4 (2%)	7 13
1	D	242/253 (96%)	223 (92%)	18 (7%)	1 (0%)	30 49
1	E	164/253 (65%)	157 (96%)	6 (4%)	1 (1%)	21 38
1	F	251/253 (99%)	226 (90%)	19 (8%)	6 (2%)	4 8

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	G	164/253 (65%)	159 (97%)	5 (3%)	0	100	100
1	H	248/253 (98%)	233 (94%)	13 (5%)	2 (1%)	16	31
All	All	1791/2024 (88%)	1652 (92%)	117 (6%)	22 (1%)	10	20

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	31	ARG
1	C	129	ASP
1	C	178	ALA
1	A	62	SER
1	B	179	ASP
1	C	132	ALA
1	D	55	ARG
1	F	223	SER
1	H	199	GLY
1	A	61	LYS
1	B	39	LEU
1	F	199	GLY
1	B	13	ALA
1	E	130	ARG
1	F	-9	SER
1	F	72	ALA
1	A	54	ASN
1	F	221	SER
1	F	10	ALA
1	H	222	VAL
1	A	66	VAL
1	C	103	VAL

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	197/198 (100%)	181 (92%)	16 (8%)	11	23

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	199/198 (100%)	177 (89%)	22 (11%)	6	12
1	C	197/198 (100%)	181 (92%)	16 (8%)	11	23
1	D	199/198 (100%)	183 (92%)	16 (8%)	11	24
1	E	134/198 (68%)	119 (89%)	15 (11%)	6	12
1	F	206/198 (104%)	185 (90%)	21 (10%)	7	15
1	G	134/198 (68%)	120 (90%)	14 (10%)	7	14
1	H	204/198 (103%)	192 (94%)	12 (6%)	18	37
All	All	1470/1584 (93%)	1338 (91%)	132 (9%)	9	19

All (132) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	LEU
1	A	15	LEU
1	A	20	GLU
1	A	31	ARG
1	A	34	LYS
1	A	79	GLN
1	A	83	ASN
1	A	92	LEU
1	A	93	ASP
1	A	126	VAL
1	A	154	VAL
1	A	177	GLU
1	A	194	VAL
1	A	222	VAL
1	A	225	LEU
1	A	235	LEU
1	B	8	ILE
1	B	12	GLN
1	B	14	LEU
1	B	28	LEU
1	B	55	ARG
1	B	58	LEU
1	B	63	ASP
1	B	83	ASN
1	B	135	ASN
1	B	143	CYS
1	B	147	GLU
1	B	154	VAL

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Mol	Chain	Res	Type
1	B	168	ASN
1	B	179	ASP
1	B	182	LEU
1	B	187	MSE
1	B	200	GLU
1	B	218	MSE
1	B	221	SER
1	B	225	LEU
1	B	226	ASN
1	B	235	LEU
1	C	4	MSE
1	C	15	LEU
1	C	20	GLU
1	C	71	ILE
1	C	73	ARG
1	C	84	ASP
1	C	92	LEU
1	C	93	ASP
1	C	130	ARG
1	C	134	LEU
1	C	143	CYS
1	C	166	GLU
1	C	172	VAL
1	C	190	ARG
1	C	224	SER
1	C	235	LEU
1	D	6	TYR
1	D	15	LEU
1	D	33	ASP
1	D	58	LEU
1	D	92	LEU
1	D	147	GLU
1	D	154	VAL
1	D	156	ASN
1	D	162	ARG
1	D	166	GLU
1	D	177	GLU
1	D	187	MSE
1	D	205	LEU
1	D	226	ASN
1	D	227	VAL
1	D	235	LEU

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Mol	Chain	Res	Type
1	E	79	GLN
1	E	83	ASN
1	E	92	LEU
1	E	94	GLN
1	E	97	LEU
1	E	126	VAL
1	E	143	CYS
1	E	147	GLU
1	E	166	GLU
1	E	177	GLU
1	E	191	LEU
1	E	205	LEU
1	E	225	LEU
1	E	235	LEU
1	E	243	SER
1	F	4	MSE
1	F	5	ILE
1	F	12	GLN
1	F	20	GLU
1	F	28	LEU
1	F	39	LEU
1	F	41	HIS
1	F	58	LEU
1	F	63	ASP
1	F	71	ILE
1	F	73	ARG
1	F	79	GLN
1	F	92	LEU
1	F	148	SER
1	F	151	LEU
1	F	154	VAL
1	F	193	LEU
1	F	225	LEU
1	F	227	VAL
1	F	235	LEU
1	F	242	ARG
1	G	79	GLN
1	G	81	GLN
1	G	93	ASP
1	G	97	LEU
1	G	130	ARG
1	G	133	GLN

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Mol	Chain	Res	Type
1	G	134	LEU
1	G	143	CYS
1	G	154	VAL
1	G	167	GLU
1	G	205	LEU
1	G	225	LEU
1	G	235	LEU
1	G	242	ARG
1	H	-6	VAL
1	H	28	LEU
1	H	63	ASP
1	H	92	LEU
1	H	143	CYS
1	H	154	VAL
1	H	162	ARG
1	H	198	GLU
1	H	200	GLU
1	H	225	LEU
1	H	235	LEU
1	H	240	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	GLN
1	A	79	GLN
1	A	94	GLN
1	A	156	ASN
1	B	79	GLN
1	B	81	GLN
1	B	94	GLN
1	B	135	ASN
1	C	9	HIS
1	C	23	GLN
1	C	46	GLN
1	C	79	GLN
1	C	81	GLN
1	C	94	GLN
1	C	133	GLN
1	C	156	ASN
1	D	81	GLN
1	D	133	GLN

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Mol	Chain	Res	Type
1	D	165	GLN
1	E	79	GLN
1	E	94	GLN
1	E	133	GLN
1	F	12	GLN
1	F	46	GLN
1	F	51	GLN
1	F	54	ASN
1	F	94	GLN
1	F	107	HIS
1	F	133	GLN
1	G	94	GLN
1	H	9	HIS
1	H	51	GLN
1	H	79	GLN
1	H	107	HIS
1	H	133	GLN
1	H	165	GLN

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 ⓘ

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	235/253 (92%)	0.68	33 (14%) 6 5	23, 53, 106, 116	1 (0%)
1	B	236/253 (93%)	0.92	33 (13%) 6 5	27, 55, 74, 116	0
1	C	235/253 (92%)	0.48	15 (6%) 25 22	21, 53, 91, 106	0
1	D	236/253 (93%)	0.51	21 (8%) 15 13	20, 45, 82, 96	0
1	E	160/253 (63%)	-0.03	0 100 100	21, 40, 63, 83	0
1	F	245/253 (96%)	0.77	56 (22%) 2 1	19, 42, 118, 126	0
1	G	160/253 (63%)	0.27	1 (0%) 85 83	26, 46, 69, 86	0
1	H	242/253 (95%)	0.35	26 (10%) 11 9	19, 43, 97, 115	0
All	All	1749/2024 (86%)	0.53	185 (10%) 11 9	19, 47, 96, 126	1 (0%)

All (185) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	28	LEU	5.6
1	F	39	LEU	5.4
1	F	10	ALA	5.3
1	F	27	ILE	5.2
1	B	18	ALA	5.1
1	A	36	LEU	4.9
1	F	69	GLY	4.8
1	F	53	ALA	4.7
1	F	36	LEU	4.7
1	F	64	GLY	4.7
1	F	28	LEU	4.6
1	B	25	VAL	4.4
1	F	58	LEU	4.3
1	F	12	GLN	4.3
1	F	57	TYR	4.2
1	F	71	ILE	4.2

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Mol	Chain	Res	Type	RSRZ
1	F	51	GLN	4.1
1	H	-6	VAL	4.1
1	F	52	LEU	4.1
1	B	69	GLY	4.0
1	F	224	SER	4.0
1	H	71	ILE	3.9
1	F	5	ILE	3.9
1	D	27	ILE	3.9
1	A	53	ALA	3.8
1	F	6	TYR	3.8
1	A	30	GLY	3.8
1	F	70	ILE	3.7
1	F	31	ARG	3.7
1	F	-5	PRO	3.7
1	F	26	PHE	3.6
1	F	25	VAL	3.6
1	F	-6	VAL	3.6
1	F	34	LYS	3.6
1	F	66	VAL	3.6
1	H	57	TYR	3.5
1	A	13	ALA	3.5
1	B	39	LEU	3.5
1	F	50	ILE	3.5
1	B	10	ALA	3.4
1	A	66	VAL	3.4
1	B	49	VAL	3.4
1	F	56	GLN	3.4
1	B	36	LEU	3.4
1	H	64	GLY	3.4
1	F	8	ILE	3.4
1	B	5	ILE	3.3
1	B	72	ALA	3.3
1	B	31	ARG	3.3
1	A	129	ASP	3.3
1	D	54	ASN	3.3
1	A	17	ARG	3.2
1	C	154	VAL	3.2
1	A	10	ALA	3.2
1	F	33	ASP	3.2
1	F	67	HIS	3.2
1	D	29	LYS	3.2
1	B	40	ILE	3.2

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Mol	Chain	Res	Type	RSRZ
1	C	17	ARG	3.2
1	F	37	LEU	3.2
1	F	54	ASN	3.2
1	B	-1	HIS	3.1
1	D	58	LEU	3.1
1	A	8	ILE	3.1
1	B	8	ILE	3.1
1	A	5	ILE	3.1
1	H	36	LEU	3.0
1	H	27	ILE	3.0
1	H	35	ARG	3.0
1	F	11	VAL	3.0
1	B	9	HIS	3.0
1	D	72	ALA	3.0
1	F	-7	LEU	3.0
1	A	38	PRO	2.9
1	B	17	ARG	2.9
1	B	19	PRO	2.9
1	A	37	LEU	2.9
1	A	6	TYR	2.9
1	F	7	GLY	2.9
1	H	-3	GLY	2.9
1	B	43	LEU	2.9
1	D	28	LEU	2.9
1	A	27	ILE	2.8
1	A	65	ALA	2.8
1	D	37	LEU	2.8
1	F	30	GLY	2.8
1	F	18	ALA	2.8
1	F	9	HIS	2.8
1	H	40	ILE	2.8
1	H	28	LEU	2.8
1	F	35	ARG	2.8
1	A	2	SER	2.8
1	H	58	LEU	2.8
1	D	33	ASP	2.8
1	A	220	GLY	2.7
1	A	61	LYS	2.7
1	H	33	ASP	2.7
1	H	26	PHE	2.7
1	B	11	VAL	2.7
1	F	65	ALA	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	71	ILE	2.7
1	F	29	LYS	2.7
1	A	52	LEU	2.7
1	C	130	ARG	2.7
1	A	222	VAL	2.6
1	A	57	TYR	2.6
1	F	68	GLN	2.6
1	F	-3	GLY	2.6
1	B	38	PRO	2.6
1	F	24	GLU	2.6
1	C	31	ARG	2.6
1	D	57	TYR	2.6
1	B	22	PHE	2.6
1	C	36	LEU	2.6
1	A	11	VAL	2.6
1	H	51	GLN	2.6
1	A	18	ALA	2.6
1	F	14	LEU	2.5
1	D	61	LYS	2.5
1	D	93	ASP	2.5
1	F	63	ASP	2.5
1	H	50	ILE	2.5
1	D	41	HIS	2.5
1	H	34	LYS	2.5
1	A	70	ILE	2.5
1	C	49	VAL	2.5
1	D	-1	HIS	2.5
1	D	38	PRO	2.5
1	B	7	GLY	2.5
1	C	199	GLY	2.5
1	D	53	ALA	2.5
1	D	69	GLY	2.4
1	F	59	ASP	2.4
1	H	6	TYR	2.4
1	H	224	SER	2.4
1	H	29	LYS	2.4
1	A	26	PHE	2.4
1	C	220	GLY	2.4
1	F	32	GLU	2.4
1	A	47	GLY	2.3
1	A	199	GLY	2.3
1	A	58	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	93	ASP	2.3
1	A	54	ASN	2.3
1	C	60	GLU	2.3
1	F	49	VAL	2.3
1	G	136	ALA	2.3
1	F	-2	SER	2.3
1	F	200	GLU	2.3
1	D	222	VAL	2.3
1	A	73	ARG	2.3
1	B	68	GLN	2.2
1	D	30	GLY	2.2
1	B	14	LEU	2.2
1	B	70	ILE	2.2
1	C	50	ILE	2.2
1	H	59	ASP	2.2
1	D	60	GLU	2.2
1	A	74	VAL	2.2
1	C	51	GLN	2.2
1	B	15	LEU	2.2
1	F	2	SER	2.2
1	H	25	VAL	2.2
1	H	70	ILE	2.2
1	H	39	LEU	2.1
1	F	-10	SER	2.1
1	F	38	PRO	2.1
1	H	41	HIS	2.1
1	C	2	SER	2.1
1	F	220	GLY	2.1
1	C	55	ARG	2.1
1	D	55	ARG	2.1
1	B	178	ALA	2.1
1	H	69	GLY	2.1
1	C	34	LYS	2.1
1	B	129	ASP	2.1
1	B	243	SER	2.1
1	B	26	PHE	2.0
1	D	66	VAL	2.0
1	B	78	ARG	2.0
1	A	224	SER	2.0
1	B	2	SER	2.0
1	H	56	GLN	2.0
1	C	66	VAL	2.0

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Mol	Chain	Res	Type	RSRZ
1	F	61	LYS	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.