



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:21 PM UTC

PDB ID : 3GKU / pdb_00003gku
Title : Crystal structure of a probable RNA-binding protein from Clostridium symbiosum ATCC 14940
Authors : Tan, K.; Keigher, L.; Jedrzejczak, R.; Babnigg, G.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2009-03-11
Resolution : 2.95 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

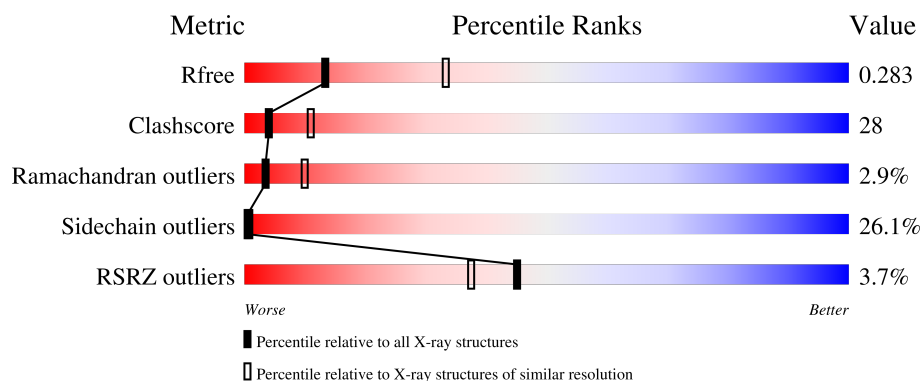
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.95 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	1130 (2.98-2.94)
Clashscore	190562	1157 (2.98-2.94)
Ramachandran outliers	187476	1101 (2.98-2.94)
Sidechain outliers	187428	1101 (2.98-2.94)
RSRZ outliers	180081	1130 (2.98-2.94)

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	225	40%40%11%8%
1	B	225	4% 41%31%13%14%
1	C	225	4% 24%29%7%40%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 4131 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 Probable RNA-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	206	Total	C	N	O	Se	0	0	0
			1636	1020	287	323	6			
1	B	193	Total	C	N	O	Se	0	0	0
			1500	934	263	297	6			
1	C	134	Total	C	N	O	Se	0	0	0
			995	619	163	208	5			

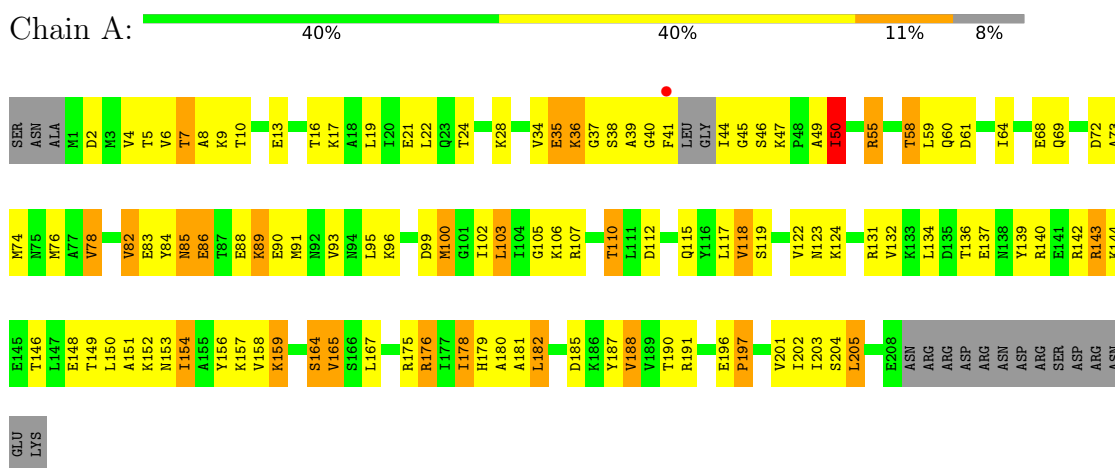
There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	PDB 3GKU
A	-1	ASN	-	expression tag	PDB 3GKU
A	0	ALA	-	expression tag	PDB 3GKU
B	-2	SER	-	expression tag	PDB 3GKU
B	-1	ASN	-	expression tag	PDB 3GKU
B	0	ALA	-	expression tag	PDB 3GKU
C	-2	SER	-	expression tag	PDB 3GKU
C	-1	ASN	-	expression tag	PDB 3GKU
C	0	ALA	-	expression tag	PDB 3GKU

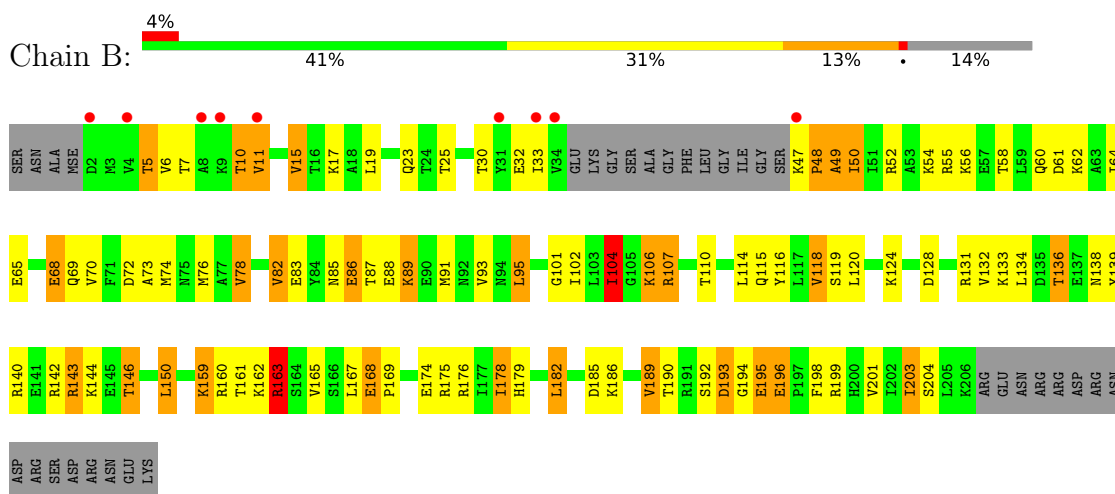
3 Residue-property plots

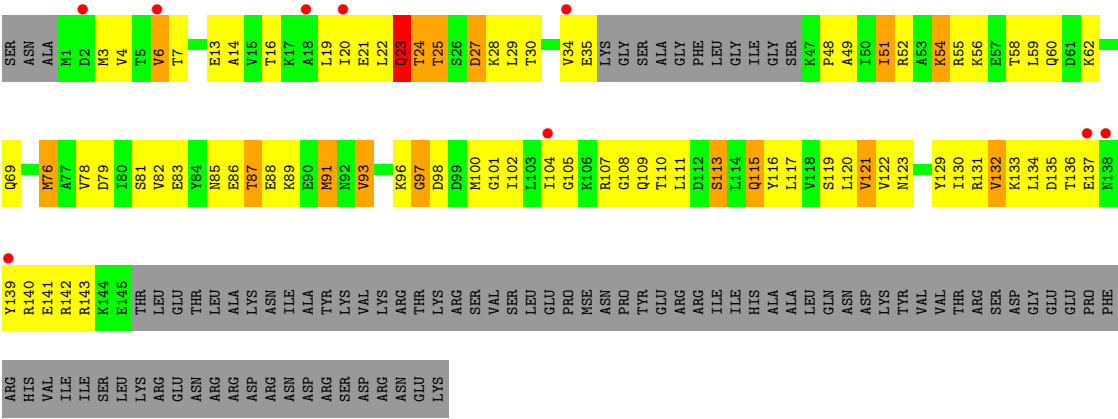
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Probable RNA-binding protein



• Molecule 1: Probable RNA-binding protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	126.99Å 126.99Å 106.19Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	40.72 – 2.95 40.72 – 2.95	Depositor EDS
% Data completeness (in resolution range)	99.8 (40.72-2.95) 99.8 (40.72-2.95)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.43 (at 2.95Å)	Xtriage
Refinement program	REFMAC 5.5.0054	Depositor
R, R_{free}	0.220 , 0.287 0.219 , 0.283	Depositor DCC
R_{free} test set	1084 reflections (5.13%)	wwPDB-VP
Wilson B-factor (Å ²)	59.8	Xtriage
Anisotropy	0.104	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 55.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.035 for -h,-k,l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	4131	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.85% 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	1.30	4/1645 (0.2%)	1.28	5/2200 (0.2%)
1	B	1.08	0/1508	1.17	5/2027 (0.2%)
1	C	0.91	0/994	1.12	4/1337 (0.3%)
All	All	1.14	4/4147 (0.1%)	1.21	14/5564 (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	C	0	1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	78	VAL	CA-CB	6.20	1.63	1.54
1	A	180	ALA	CA-CB	-5.32	1.44	1.53
1	A	149	THR	CA-CB	-5.29	1.44	1.53
1	A	202	ILE	CA-CB	-5.06	1.48	1.54

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	97	GLY	N-CA-C	8.41	129.37	115.66
1	C	135	ASP	N-CA-C	7.54	120.63	109.24
1	B	136	THR	CB-CA-C	-6.64	99.40	110.29
1	A	105	GLY	N-CA-C	-6.61	104.31	112.77
1	A	72	ASP	N-CA-C	-6.54	103.36	112.45
1	A	123	ASN	N-CA-C	6.54	119.41	111.82
1	B	72	ASP	N-CA-C	-6.50	104.27	111.36
1	C	123	ASN	N-CA-C	6.48	119.33	111.82
1	A	165	VAL	N-CA-C	6.47	117.17	108.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	193	ASP	N-CA-C	6.02	117.29	108.14
1	A	50	ILE	CB-CA-C	-5.72	101.97	110.33
1	C	108	GLY	N-CA-C	-5.41	107.56	114.37
1	B	50	ILE	N-CA-C	5.25	116.94	108.85
1	B	19	LEU	N-CA-C	-5.19	105.62	111.28

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	24	THR	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	1636	0	1673	89	0
1	B	1500	0	1483	68	0
1	C	995	0	949	77	0
All	All	4131	0	4105	230	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 28.

All (230) 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:103:LEU:HD23	1:A:103:LEU:H	1.02	1.13
1:C:100:MSE:HE2	1:C:136:THR:HG22	1.24	1.12
1:A:110:THR:HG22	1:B:116:TYR:HE2	1.08	1.10
1:C:97:GLY:HA2	1:C:100:MSE:HG2	1.29	1.10
1:C:100:MSE:HE2	1:C:136:THR:CG2	1.86	1.05
1:A:103:LEU:H	1:A:103:LEU:CD2	1.74	0.99
1:A:110:THR:HG22	1:B:116:TYR:CE2	1.97	0.98
1:B:95:LEU:H	1:B:95:LEU:CD2	1.75	0.98

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:106:LYS:O	1:A:110:THR:HG23	1.63	0.98
1:C:100:MSE:CE	1:C:136:THR:CG2	2.43	0.96
1:C:82:VAL:HG21	1:C:91:MSE:HE2	1.48	0.96
1:C:101:GLY:O	1:C:104:ILE:HB	1.65	0.95
1:A:69:GLN:HG2	1:B:73:ALA:HB2	1.51	0.92
1:B:95:LEU:H	1:B:95:LEU:HD22	1.36	0.89
1:C:97:GLY:HA2	1:C:100:MSE:CG	2.02	0.89
1:A:103:LEU:HD23	1:A:103:LEU:N	1.84	0.89
1:B:78:VAL:HG21	1:B:95:LEU:HB2	1.54	0.87
1:A:58:THR:HG22	1:A:61:ASP:H	1.40	0.87
1:B:182:LEU:HD23	1:B:203:ILE:HD13	1.55	0.86
1:B:95:LEU:HD22	1:B:95:LEU:N	1.92	0.84
1:C:100:MSE:CE	1:C:136:THR:HG21	2.08	0.84
1:A:74:MSE:HE2	1:A:76:MSE:SE	2.28	0.83
1:B:82:VAL:HB	1:B:93:VAL:HG22	1.61	0.83
1:B:95:LEU:CD2	1:B:95:LEU:N	2.36	0.82
1:C:97:GLY:CA	1:C:100:MSE:HG2	2.07	0.82
1:A:73:ALA:HB2	1:B:69:GLN:HG2	1.62	0.81
1:C:76:MSE:HE1	1:C:102:ILE:HG21	1.63	0.80
1:B:182:LEU:HD23	1:B:203:ILE:CD1	2.13	0.79
1:A:154:ILE:HD13	1:A:165:VAL:HG11	1.65	0.79
1:C:60:GLN:HA	1:C:91:MSE:HE1	1.66	0.78
1:C:23:GLN:OE1	1:C:23:GLN:HA	1.84	0.78
1:C:119:SER:HA	1:C:132:VAL:HG21	1.65	0.78
1:C:100:MSE:CE	1:C:136:THR:HG22	2.04	0.76
1:A:10:THR:OG1	1:A:13:GLU:HG3	1.85	0.76
1:C:100:MSE:HE3	1:C:136:THR:HG21	1.65	0.76
1:A:44:ILE:O	1:A:46:SER:N	2.18	0.75
1:A:136:THR:O	1:A:137:GLU:C	2.29	0.75
1:B:95:LEU:H	1:B:95:LEU:HD23	1.52	0.74
1:B:78:VAL:CG2	1:B:95:LEU:HB2	2.18	0.74
1:A:39:ALA:HA	1:A:46:SER:HA	1.70	0.74
1:C:58:THR:HG21	1:C:60:GLN:HG2	1.68	0.73
1:C:119:SER:HB2	1:C:132:VAL:CG2	2.18	0.73
1:A:175:ARG:HG2	1:A:201:VAL:HG23	1.69	0.72
1:A:93:VAL:HG13	1:A:134:LEU:HD12	1.74	0.70
1:C:51:ILE:HG12	1:C:52:ARG:N	2.05	0.70
1:C:35:GLU:HG2	1:C:49:ALA:HA	1.74	0.69
1:C:137:GLU:HA	1:C:137:GLU:OE1	1.93	0.69
1:C:117:LEU:O	1:C:121:VAL:HG23	1.93	0.69
1:C:24:THR:HB	1:C:55:ARG:HH11	1.57	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:154:ILE:HD13	1:A:165:VAL:CG1	2.23	0.68
1:C:27:ASP:OD1	1:C:27:ASP:N	2.23	0.68
1:A:142:ARG:O	1:A:146:THR:HG23	1.93	0.68
1:C:119:SER:HB2	1:C:132:VAL:HG23	1.76	0.68
1:A:58:THR:O	1:A:61:ASP:HB2	1.94	0.67
1:C:58:THR:HG22	1:C:60:GLN:H	1.58	0.67
1:A:196:GLU:HG3	1:A:197:PRO:HA	1.77	0.67
1:B:179:HIS:ND1	1:B:190:THR:OG1	2.24	0.67
1:C:22:LEU:CB	1:C:55:ARG:HH12	2.09	0.66
1:C:25:THR:HB	1:C:27:ASP:OD1	1.95	0.65
1:A:91:MSE:HB3	1:A:132:VAL:HG12	1.79	0.65
1:A:93:VAL:HG13	1:A:134:LEU:CD1	2.27	0.65
1:C:58:THR:HG21	1:C:60:GLN:CG	2.27	0.64
1:B:182:LEU:CD2	1:B:203:ILE:CD1	2.75	0.64
1:A:139:TYR:CZ	1:A:143:ARG:HG2	2.34	0.63
1:A:34:VAL:HB	1:A:50:ILE:HB	1.81	0.63
1:A:85:ASN:C	1:A:85:ASN:HD22	2.07	0.62
1:C:119:SER:CA	1:C:132:VAL:HG21	2.30	0.61
1:A:144:LYS:HE2	1:A:148:GLU:OE2	2.01	0.61
1:A:164:SER:OG	1:A:204:SER:HB3	2.01	0.61
1:A:112:ASP:OD1	1:A:140:ARG:NH2	2.34	0.60
1:C:141:GLU:C	1:C:143:ARG:H	2.08	0.60
1:A:99:ASP:O	1:A:102:ILE:HG12	2.01	0.60
1:B:193:ASP:OD2	1:B:194:GLY:N	2.34	0.60
1:A:24:THR:OG1	1:A:28:LYS:HB2	2.01	0.60
1:A:84:TYR:CE2	1:A:86:GLU:HG2	2.37	0.59
1:A:205:LEU:CD2	1:A:205:LEU:H	2.15	0.59
1:C:119:SER:CB	1:C:132:VAL:HG21	2.31	0.59
1:C:22:LEU:O	1:C:24:THR:N	2.30	0.59
1:B:65:GLU:OE1	1:B:65:GLU:HA	2.01	0.59
1:B:58:THR:HG22	1:B:60:GLN:N	2.18	0.58
1:B:58:THR:HG22	1:B:60:GLN:H	1.68	0.58
1:A:205:LEU:H	1:A:205:LEU:HD22	1.67	0.57
1:C:58:THR:HG22	1:C:59:LEU:N	2.21	0.56
1:C:129:TYR:HE2	1:C:131:ARG:HD2	1.69	0.56
1:B:64:ILE:O	1:B:68:GLU:HB2	2.05	0.56
1:B:70:VAL:O	1:B:74:MSE:HG3	2.06	0.56
1:B:32:GLU:HB2	1:B:52:ARG:HB2	1.88	0.55
1:A:100:MSE:HA	1:A:100:MSE:CE	2.37	0.55
1:A:103:LEU:O	1:A:107:ARG:HG3	2.06	0.55
1:C:79:ASP:HB2	1:C:96:LYS:CB	2.36	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:93:VAL:CG1	1:B:134:LEU:CD1	2.85	0.55
1:B:161:THR:C	1:B:163:ARG:H	2.15	0.55
1:A:38:SER:O	1:A:46:SER:HA	2.06	0.55
1:C:107:ARG:HB2	1:C:107:ARG:HH11	1.72	0.55
1:B:142:ARG:O	1:B:146:THR:HG23	2.07	0.54
1:C:119:SER:CB	1:C:132:VAL:CG2	2.86	0.54
1:B:85:ASN:ND2	1:B:88:GLU:HG2	2.21	0.54
1:B:101:GLY:O	1:B:104:ILE:HG12	2.08	0.54
1:C:119:SER:HB2	1:C:132:VAL:HG21	1.86	0.54
1:B:169:PRO:HB3	1:B:198:PHE:HA	1.91	0.53
1:C:115:GLN:HG3	1:C:134:LEU:HB2	1.90	0.53
1:B:86:GLU:O	1:B:89:LYS:HD3	2.09	0.53
1:A:37:GLY:CA	1:A:47:LYS:O	2.58	0.52
1:C:141:GLU:C	1:C:143:ARG:N	2.68	0.52
1:C:3:MSE:HG3	1:C:54:LYS:HG2	1.92	0.52
1:A:7:THR:O	1:A:8:ALA:HB2	2.09	0.51
1:A:58:THR:HG21	1:A:60:GLN:HG2	1.92	0.51
1:C:3:MSE:HE2	1:C:54:LYS:HZ3	1.74	0.51
1:B:150:LEU:HD23	1:B:150:LEU:O	2.10	0.51
1:C:6:VAL:HG13	1:C:51:ILE:HD11	1.92	0.51
1:B:95:LEU:N	1:B:95:LEU:HD23	2.17	0.51
1:C:3:MSE:HE2	1:C:54:LYS:NZ	2.26	0.51
1:A:85:ASN:O	1:A:89:LYS:N	2.42	0.50
1:A:102:ILE:HG13	1:A:103:LEU:HD22	1.92	0.50
1:C:16:THR:O	1:C:20:ILE:HG13	2.12	0.50
1:A:88:GLU:OE2	1:A:88:GLU:HA	2.07	0.50
1:B:58:THR:CG2	1:B:60:GLN:OE1	2.60	0.50
1:C:115:GLN:HG3	1:C:134:LEU:CB	2.42	0.49
1:B:58:THR:HB	1:B:61:ASP:CG	2.37	0.49
1:B:159:LYS:HD2	1:B:159:LYS:C	2.37	0.49
1:A:154:ILE:CD1	1:A:165:VAL:HB	2.42	0.49
1:A:58:THR:HG23	1:A:60:GLN:OE1	2.13	0.49
1:B:139:TYR:O	1:B:140:ARG:C	2.54	0.49
1:A:40:GLY:C	1:A:44:ILE:HA	2.38	0.49
1:A:119:SER:HB2	1:A:132:VAL:CG2	2.42	0.49
1:A:167:LEU:N	1:A:167:LEU:CD1	2.75	0.49
1:B:5:THR:HG23	1:B:52:ARG:HE	1.77	0.49
1:B:5:THR:HG23	1:B:52:ARG:NE	2.28	0.49
1:A:158:VAL:O	1:A:159:LYS:C	2.55	0.48
1:C:89:LYS:HG2	1:C:130:ILE:HD12	1.95	0.48
1:C:105:GLY:HA2	1:C:110:THR:HG23	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:107:ARG:HB2	1:C:107:ARG:NH1	2.27	0.48
1:A:7:THR:O	1:A:17:LYS:HE2	2.13	0.48
1:C:129:TYR:CE2	1:C:131:ARG:HD2	2.48	0.48
1:A:102:ILE:HG13	1:A:103:LEU:N	2.29	0.48
1:C:116:TYR:C	1:C:116:TYR:CD2	2.92	0.48
1:A:178:ILE:O	1:A:179:HIS:C	2.57	0.48
1:A:60:GLN:O	1:A:64:ILE:HG13	2.14	0.48
1:B:139:TYR:CZ	1:B:143:ARG:HG2	2.49	0.48
1:C:51:ILE:CG1	1:C:52:ARG:N	2.76	0.48
1:B:93:VAL:HG12	1:B:134:LEU:CD1	2.43	0.48
1:C:14:ALA:HB3	1:C:51:ILE:HG23	1.96	0.48
1:A:100:MSE:HA	1:A:100:MSE:HE3	1.96	0.47
1:B:150:LEU:C	1:B:150:LEU:CD2	2.87	0.47
1:C:82:VAL:HG22	1:C:83:GLU:N	2.29	0.47
1:C:82:VAL:HA	1:C:93:VAL:HA	1.96	0.47
1:A:90:GLU:HB3	1:A:131:ARG:HB3	1.97	0.47
1:B:138:ASN:O	1:B:139:TYR:C	2.57	0.47
1:C:85:ASN:ND2	1:C:88:GLU:HB2	2.29	0.47
1:B:11:VAL:O	1:B:15:VAL:N	2.48	0.46
1:A:41:PHE:HA	1:A:44:ILE:HA	1.98	0.46
1:A:103:LEU:CD2	1:A:103:LEU:N	2.52	0.46
1:A:176:ARG:O	1:A:176:ARG:HG3	2.07	0.46
1:A:10:THR:HG1	1:A:13:GLU:HG3	1.80	0.46
1:A:55:ARG:HE	1:A:55:ARG:HB2	1.60	0.46
1:A:35:GLU:O	1:A:49:ALA:HA	2.16	0.46
1:A:151:ALA:HB1	1:A:182:LEU:HD13	1.97	0.46
1:A:178:ILE:O	1:A:181:ALA:N	2.48	0.46
1:C:109:GLN:O	1:C:113:SER:HB2	2.16	0.46
1:B:65:GLU:OE1	1:B:65:GLU:CA	2.64	0.46
1:C:86:GLU:O	1:C:87:THR:C	2.56	0.46
1:A:38:SER:O	1:A:47:LYS:N	2.47	0.46
1:A:58:THR:CG2	1:A:60:GLN:HG2	2.46	0.46
1:B:161:THR:C	1:B:163:ARG:N	2.74	0.46
1:A:58:THR:HB	1:A:61:ASP:CG	2.41	0.45
1:B:10:THR:HG23	1:B:11:VAL:H	1.81	0.45
1:A:37:GLY:HA3	1:A:47:LYS:O	2.16	0.45
1:B:182:LEU:HD12	1:B:182:LEU:HA	1.81	0.45
1:B:195:GLU:O	1:B:196:GLU:C	2.59	0.45
1:C:101:GLY:HA2	1:C:104:ILE:HD12	1.99	0.45
1:A:88:GLU:O	1:A:89:LYS:CB	2.63	0.44
1:A:95:LEU:O	1:A:136:THR:HG23	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:106:LYS:HB2	1:B:106:LYS:HE2	1.74	0.44
1:A:36:LYS:H	1:A:36:LYS:HG2	1.62	0.44
1:C:97:GLY:O	1:C:98:ASP:C	2.60	0.44
1:A:82:VAL:HG23	1:A:83:GLU:N	2.33	0.44
1:A:6:VAL:HG12	1:A:7:THR:N	2.31	0.43
1:A:76:MSE:HE2	1:A:102:ILE:HD11	2.00	0.43
1:B:76:MSE:HE2	1:B:102:ILE:HB	1.98	0.43
1:A:85:ASN:C	1:A:85:ASN:ND2	2.76	0.43
1:C:58:THR:HB	1:C:60:GLN:HG3	2.00	0.43
1:B:93:VAL:CG1	1:B:134:LEU:HD12	2.47	0.43
1:B:7:THR:HG23	1:B:49:ALA:O	2.18	0.43
1:C:97:GLY:HA2	1:C:100:MSE:SE	2.69	0.43
1:A:117:LEU:O	1:A:118:VAL:C	2.61	0.43
1:A:182:LEU:HD23	1:A:190:THR:CG2	2.48	0.42
1:B:139:TYR:HA	1:B:142:ARG:NH2	2.34	0.42
1:A:153:ASN:O	1:A:156:TYR:HB3	2.19	0.42
1:C:82:VAL:CG2	1:C:83:GLU:N	2.82	0.42
1:A:158:VAL:HG21	1:A:203:ILE:HG22	2.01	0.42
1:C:6:VAL:HG13	1:C:51:ILE:CD1	2.50	0.42
1:A:159:LYS:HD2	1:A:187:TYR:HB2	2.02	0.42
1:A:205:LEU:CD2	1:A:205:LEU:N	2.82	0.42
1:A:85:ASN:CG	1:A:88:GLU:HB2	2.45	0.42
1:B:85:ASN:ND2	1:B:88:GLU:CG	2.83	0.41
1:C:22:LEU:C	1:C:24:THR:N	2.78	0.41
1:A:190:THR:HG22	1:A:203:ILE:HG12	2.01	0.41
1:A:59:LEU:HD22	1:A:122:VAL:HG13	2.02	0.41
1:C:104:ILE:HG22	1:C:105:GLY:N	2.35	0.41
1:A:115:GLN:OE1	1:A:134:LEU:N	2.47	0.41
1:B:175:ARG:HG2	1:B:201:VAL:HG23	2.01	0.41
1:C:60:GLN:CA	1:C:91:MSE:HE1	2.41	0.41
1:A:85:ASN:ND2	1:A:88:GLU:H	2.19	0.41
1:B:60:GLN:O	1:B:61:ASP:C	2.63	0.41
1:B:118:VAL:CG1	1:B:119:SER:N	2.84	0.41
1:A:58:THR:HB	1:A:61:ASP:OD2	2.21	0.41
1:A:185:ASP:OD2	1:A:188:VAL:N	2.52	0.41
1:B:48:PRO:O	1:B:49:ALA:HB2	2.21	0.41
1:B:54:LYS:HE2	1:B:56:LYS:HA	2.03	0.41
1:B:91:MSE:HE3	1:B:91:MSE:HB2	1.93	0.41
1:B:189:VAL:HG13	1:B:204:SER:O	2.21	0.41
1:C:29:LEU:HG	1:C:30:THR:N	2.36	0.41
1:C:139:TYR:O	1:C:140:ARG:C	2.63	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:141:GLU:O	1:C:143:ARG:N	2.54	0.41
1:B:114:LEU:O	1:B:115:GLN:C	2.61	0.41
1:B:167:LEU:HD13	1:B:201:VAL:HB	2.02	0.41
1:C:105:GLY:HA3	1:C:110:THR:H	1.86	0.41
1:B:7:THR:HA	1:B:49:ALA:O	2.20	0.40
1:B:47:LYS:N	1:B:48:PRO:CD	2.85	0.40
1:C:25:THR:HB	1:C:27:ASP:CG	2.47	0.40
1:C:24:THR:OG1	1:C:28:LYS:CB	2.69	0.40
1:C:129:TYR:CD2	1:C:129:TYR:C	2.98	0.40
1:C:59:LEU:HD23	1:C:59:LEU:HA	1.89	0.40
1:B:168:GLU:H	1:B:168:GLU:HG3	1.39	0.40
1:C:58:THR:CG2	1:C:60:GLN:H	2.30	0.40
1:A:22:LEU:HA	1:A:22:LEU:HD23	1.79	0.40
1:A:178:ILE:HG22	1:A:179:HIS:N	2.36	0.40
1:B:93:VAL:HG12	1:B:134:LEU:HD12	2.03	0.40
1:B:178:ILE:O	1:B:179:HIS:C	2.62	0.40
1:C:121:VAL:O	1:C:122:VAL:C	2.64	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	202/225 (90%)	179 (89%)	22 (11%)	1 (0%)	24	49
1	B	189/225 (84%)	161 (85%)	17 (9%)	11 (6%)	1	2
1	C	130/225 (58%)	110 (85%)	17 (13%)	3 (2%)	5	14
All	All	521/675 (77%)	450 (86%)	56 (11%)	15 (3%)	3	10

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	45	GLY
1	B	107	ARG
1	C	23	GLN
1	C	48	PRO
1	B	17	LYS
1	B	23	GLN
1	B	163	ARG
1	B	186	LYS
1	B	162	LYS
1	B	185	ASP
1	C	142	ARG
1	B	49	ALA
1	B	48	PRO
1	B	11	VAL
1	B	104	ILE

5.3.2 Protein sidechains [i](#)

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	181/195 (93%)	142 (78%)	39 (22%)	1	2
1	B	161/195 (83%)	112 (70%)	49 (30%)	0	0
1	C	103/195 (53%)	75 (73%)	28 (27%)	0	0
All	All	445/585 (76%)	329 (74%)	116 (26%)	0	1

All (116) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ASP
1	A	4	VAL
1	A	5	THR
1	A	7	THR
1	A	9	LYS
1	A	16	THR
1	A	19	LEU
1	A	21	GLU

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Mol	Chain	Res	Type
1	A	35	GLU
1	A	36	LYS
1	A	50	ILE
1	A	55	ARG
1	A	58	THR
1	A	68	GLU
1	A	78	VAL
1	A	82	VAL
1	A	85	ASN
1	A	86	GLU
1	A	89	LYS
1	A	96	LYS
1	A	100	MSE
1	A	103	LEU
1	A	110	THR
1	A	118	VAL
1	A	124	LYS
1	A	143	ARG
1	A	150	LEU
1	A	152	LYS
1	A	154	ILE
1	A	157	LYS
1	A	159	LYS
1	A	164	SER
1	A	176	ARG
1	A	178	ILE
1	A	182	LEU
1	A	188	VAL
1	A	191	ARG
1	A	197	PRO
1	A	205	LEU
1	B	5	THR
1	B	6	VAL
1	B	10	THR
1	B	15	VAL
1	B	25	THR
1	B	30	THR
1	B	33	ILE
1	B	50	ILE
1	B	55	ARG
1	B	62	LYS
1	B	68	GLU

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Mol	Chain	Res	Type
1	B	78	VAL
1	B	82	VAL
1	B	83	GLU
1	B	86	GLU
1	B	87	THR
1	B	89	LYS
1	B	95	LEU
1	B	104	ILE
1	B	106	LYS
1	B	107	ARG
1	B	110	THR
1	B	118	VAL
1	B	120	LEU
1	B	124	LYS
1	B	128	ASP
1	B	131	ARG
1	B	132	VAL
1	B	133	LYS
1	B	136	THR
1	B	143	ARG
1	B	144	LYS
1	B	146	THR
1	B	150	LEU
1	B	159	LYS
1	B	160	ARG
1	B	163	ARG
1	B	165	VAL
1	B	168	GLU
1	B	174	GLU
1	B	176	ARG
1	B	178	ILE
1	B	182	LEU
1	B	189	VAL
1	B	192	SER
1	B	195	GLU
1	B	196	GLU
1	B	199	ARG
1	B	203	ILE
1	C	4	VAL
1	C	6	VAL
1	C	7	THR
1	C	13	GLU

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Mol	Chain	Res	Type
1	C	19	LEU
1	C	21	GLU
1	C	23	GLN
1	C	25	THR
1	C	27	ASP
1	C	34	VAL
1	C	51	ILE
1	C	54	LYS
1	C	56	LYS
1	C	62	LYS
1	C	69	GLN
1	C	76	MSE
1	C	78	VAL
1	C	81	SER
1	C	87	THR
1	C	91	MSE
1	C	93	VAL
1	C	111	LEU
1	C	113	SER
1	C	115	GLN
1	C	120	LEU
1	C	121	VAL
1	C	132	VAL
1	C	133	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	85	ASN
1	A	153	ASN
1	A	183	GLN
1	A	184	ASN
1	A	200	HIS
1	B	75	ASN
1	B	153	ASN
1	B	200	HIS
1	C	60	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no 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	199/225 (88%)	-0.32	1 (0%) 87 83	17, 32, 56, 70	0
1	B	187/225 (83%)	0.23	9 (4%) 35 30	27, 46, 82, 92	0
1	C	128/225 (56%)	0.59	9 (7%) 22 19	38, 56, 71, 76	0
All	All	514/675 (76%)	0.11	19 (3%) 45 37	17, 44, 72, 92	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	138	ASN	3.6
1	B	47	LYS	3.4
1	C	2	ASP	3.2
1	B	2	ASP	2.9
1	C	104	ILE	2.9
1	C	34	VAL	2.8
1	B	8	ALA	2.7
1	B	4	VAL	2.7
1	A	41	PHE	2.7
1	C	20	ILE	2.6
1	B	33	ILE	2.6
1	B	11	VAL	2.6
1	B	34	VAL	2.4
1	C	6	VAL	2.4
1	C	137	GLU	2.3
1	C	139	TYR	2.2
1	B	9	LYS	2.2
1	C	18	ALA	2.1
1	B	31	TYR	2.1

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.