



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 10:10 AM UTC

PDB ID : 3KIO / pdb_00003kio
Title : mouse RNase H2 complex
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Deposited on : 2009-11-02
Resolution : 2.90 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

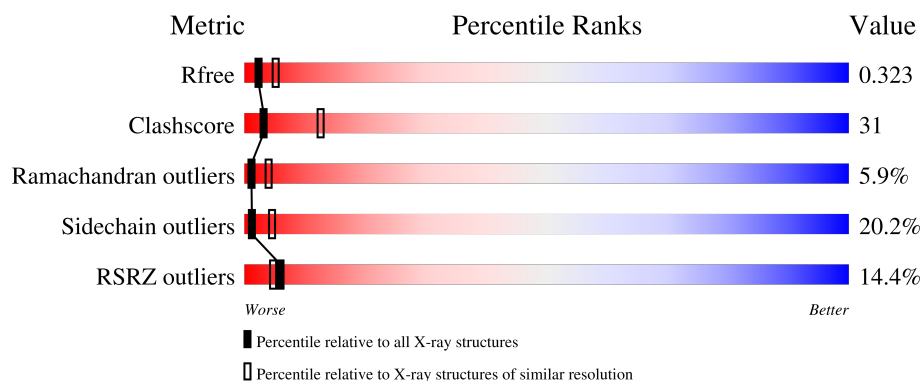
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	301	6% 39% 34% 11% 15%
2	B	332	10% 21% 15% 5% 58%
3	C	166	11% 27% 23% 14% 32%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3854 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 Ribonuclease H2 subunit A.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	256	Total	C	N	O	S	Se	0	0	0
			1986	1261	331	384	5	5			

- Molecule 2 is a protein called Ribonuclease H2 subunit B.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	140	Total	C	N	O	S	Se	0	0	0
			983	636	170	173	1	3			

There are 25 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-23	MSE	-	expression tag	UNP Q80ZV0
B	-22	GLY	-	expression tag	UNP Q80ZV0
B	-21	SER	-	expression tag	UNP Q80ZV0
B	-20	SER	-	expression tag	UNP Q80ZV0
B	-19	HIS	-	expression tag	UNP Q80ZV0
B	-18	HIS	-	expression tag	UNP Q80ZV0
B	-17	HIS	-	expression tag	UNP Q80ZV0
B	-16	HIS	-	expression tag	UNP Q80ZV0
B	-15	HIS	-	expression tag	UNP Q80ZV0
B	-14	HIS	-	expression tag	UNP Q80ZV0
B	-13	SER	-	expression tag	UNP Q80ZV0
B	-12	GLN	-	expression tag	UNP Q80ZV0
B	-11	ASP	-	expression tag	UNP Q80ZV0
B	-10	PRO	-	expression tag	UNP Q80ZV0
B	-9	ASN	-	expression tag	UNP Q80ZV0
B	-8	SER	-	expression tag	UNP Q80ZV0
B	-7	LEU	-	expression tag	UNP Q80ZV0
B	-6	GLU	-	expression tag	UNP Q80ZV0
B	-5	VAL	-	expression tag	UNP Q80ZV0
B	-4	LEU	-	expression tag	UNP Q80ZV0

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-3	PHE	-	expression tag	UNP Q80ZV0
B	-2	GLN	-	expression tag	UNP Q80ZV0
B	-1	GLY	-	expression tag	UNP Q80ZV0
B	0	PRO	-	expression tag	UNP Q80ZV0
B	239	PRO	THR	variant	UNP Q80ZV0

- Molecule 3 is a protein called Ribonuclease H2 subunit C.

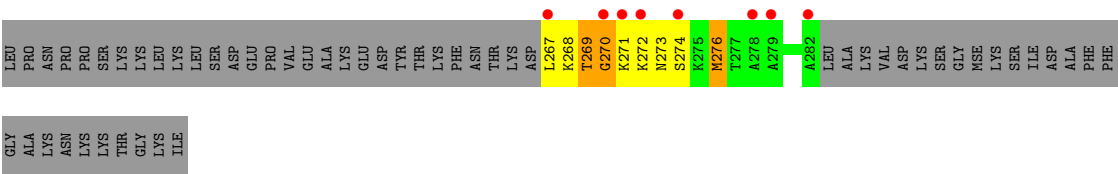
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	C	113	Total	C	N	O	S	Se	0	0	0
			825	521	149	153	1	1			

There is a discrepancy between the modelled and reference sequences:

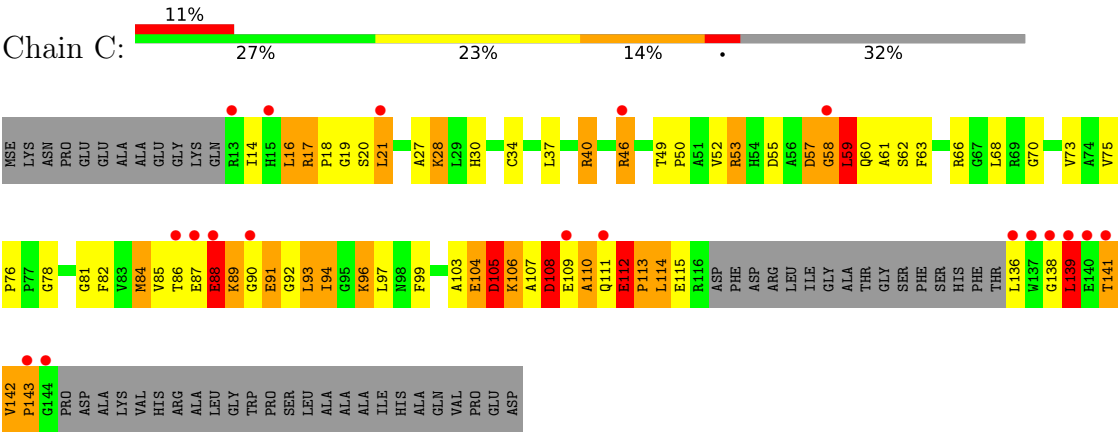
Chain	Residue	Modelled	Actual	Comment	Reference
C	9	GLU	ASP	variant	UNP Q9CQ18

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	42	Total	O	0	0
			42	42		
4	B	6	Total	O	0	0
			6	6		
4	C	12	Total	O	0	0
			12	12		



● Molecule 3: Ribonuclease H2 subunit C



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	279.29Å 40.42Å 67.82Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.62 – 2.90 24.62 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.3 (24.62-2.90) 99.1 (24.62-2.90)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	8.90 (at 2.89Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.4_113)	Depositor
R, R_{free}	0.234 , 0.305 0.250 , 0.323	Depositor DCC
R_{free} test set	925 reflections (5.20%)	wwPDB-VP
Wilson B-factor (Å ²)	49.9	Xtriage
Anisotropy	0.011	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 72.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	3854	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.20% 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	0.74	5/2025 (0.2%)	1.17	21/2751 (0.8%)
2	B	0.59	2/997 (0.2%)	0.97	5/1354 (0.4%)
3	C	0.62	0/842	1.10	5/1140 (0.4%)
All	All	0.68	7/3864 (0.2%)	1.10	31/5245 (0.6%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	B	0	1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	112	TYR	C-O	-5.46	1.17	1.23
1	A	1	MSE	SE-CE	-5.30	1.79	1.95
1	A	146	MSE	SE-CE	-5.27	1.79	1.95
2	B	55	MSE	SE-CE	-5.27	1.79	1.95
2	B	276	MSE	SE-CE	-5.26	1.79	1.95
1	A	114	LEU	CA-C	-5.09	1.46	1.52
1	A	1	MSE	CG-SE	-5.03	1.80	1.95

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	64	LYS	N-CA-C	13.42	131.86	112.94
1	A	231	PHE	CA-C-N	8.90	128.94	119.78
1	A	231	PHE	C-N-CA	8.90	128.94	119.78
1	A	64	LYS	N-CA-CB	-8.46	97.02	110.98
1	A	111	LYS	N-CA-C	8.23	122.16	110.23
1	A	111	LYS	CB-CA-C	-7.25	97.69	109.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	83	GLY	N-CA-C	-7.23	105.08	115.27
1	A	89	PHE	N-CA-CB	-7.22	100.56	110.95
1	A	39	GLY	N-CA-C	7.22	127.06	112.34
3	C	58	GLY	N-CA-C	7.10	130.00	113.18
1	A	214	ASP	CA-C-N	6.89	128.45	119.84
1	A	214	ASP	C-N-CA	6.89	128.45	119.84
3	C	40	ARG	CA-C-N	-6.89	112.55	119.92
3	C	40	ARG	C-N-CA	-6.89	112.55	119.92
1	A	46	VAL	N-CA-C	6.44	117.13	108.11
1	A	8	ARG	N-CA-C	-6.31	105.30	112.87
1	A	88	ASP	CB-CA-C	6.21	122.07	110.70
2	B	101	LEU	N-CA-C	-6.20	106.36	114.04
1	A	113	ASN	N-CA-C	6.19	119.56	110.59
1	A	231	PHE	N-CA-C	6.12	117.16	110.13
1	A	43	GLY	CA-C-N	5.88	125.79	119.85
1	A	43	GLY	C-N-CA	5.88	125.79	119.85
3	C	59	LEU	N-CA-CB	5.63	120.11	110.77
1	A	173	PHE	CA-C-N	5.62	125.47	119.28
1	A	173	PHE	C-N-CA	5.62	125.47	119.28
1	A	212	PRO	N-CA-C	-5.57	103.05	111.41
2	B	98	LEU	N-CA-C	-5.54	106.88	113.97
1	A	89	PHE	N-CA-C	5.50	122.59	113.89
2	B	92	ASP	CA-C-N	5.19	126.33	119.84
2	B	92	ASP	C-N-CA	5.19	126.33	119.84
3	C	112	GLU	N-CA-C	-5.19	100.05	109.09

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	92	ASP	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	1986	0	1927	106	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	983	0	866	62	0
3	C	825	0	795	78	0
4	A	42	0	0	3	0
4	B	6	0	0	0	0
4	C	12	0	0	1	0
All	All	3854	0	3588	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 31.

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:222:ARG:HG2	1:A:222:ARG:HH11	1.06	1.20
2:B:93:PRO:HD3	2:B:270:GLY:HA2	1.41	1.02
1:A:27:GLU:HG3	1:A:28:PRO:HD2	1.45	0.96
1:A:2:ASP:HA	1:A:5:GLU:CG	2.01	0.91
1:A:27:GLU:HG3	1:A:28:PRO:CD	2.08	0.84
3:C:58:GLY:O	3:C:59:LEU:HG	1.77	0.84
1:A:222:ARG:HG2	1:A:222:ARG:NH1	1.87	0.81
1:A:33:VAL:HG22	1:A:141:VAL:HG13	1.60	0.81
2:B:268:LYS:HG3	2:B:269:THR:H	1.46	0.80
3:C:21:LEU:H	3:C:21:LEU:HD23	1.46	0.80
1:A:109:ARG:HH22	3:C:106:LYS:HG2	1.48	0.78
1:A:194:TRP:CH2	1:A:198:GLU:HG3	2.20	0.76
1:A:2:ASP:HA	1:A:5:GLU:HG3	1.67	0.76
2:B:177:ALA:C	2:B:178:ASN:OD1	2.30	0.74
2:B:163:TRP:HA	2:B:267:LEU:HD11	1.69	0.74
1:A:109:ARG:NH2	3:C:106:LYS:HG2	2.02	0.73
2:B:162:LYS:HA	2:B:165:GLU:HB3	1.68	0.73
2:B:101:LEU:HD21	2:B:178:ASN:HD22	1.53	0.73
1:A:31:LEU:HD12	1:A:51:TYR:HB3	1.69	0.72
3:C:53:ARG:HG2	3:C:53:ARG:HH11	1.52	0.72
1:A:41:VAL:HB	1:A:239:TRP:CD1	2.25	0.72
3:C:58:GLY:HA3	3:C:70:GLY:O	1.90	0.71
1:A:147:PRO:HB3	1:A:165:VAL:HG12	1.71	0.71
2:B:93:PRO:CD	2:B:270:GLY:HA2	2.18	0.71
3:C:53:ARG:HH11	3:C:53:ARG:CG	2.04	0.70
2:B:17:PHE:HD2	3:C:84:MSE:HG3	1.56	0.70
3:C:52:VAL:C	3:C:53:ARG:HD2	2.17	0.70
1:A:1:MSE:HG2	1:A:3:LEU:HD12	1.73	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:233:GLN:HA	1:A:233:GLN:OE1	1.90	0.70
3:C:55:ASP:O	3:C:57:ASP:O	2.10	0.69
1:A:33:VAL:CG2	1:A:141:VAL:HG13	2.23	0.68
1:A:222:ARG:HH11	1:A:222:ARG:CG	1.93	0.68
2:B:93:PRO:HB2	2:B:170:GLN:NE2	2.08	0.68
1:A:156:GLN:O	1:A:159:PRO:HG3	1.94	0.67
2:B:101:LEU:HD21	2:B:178:ASN:ND2	2.08	0.67
1:A:194:TRP:CZ3	1:A:198:GLU:HG3	2.30	0.67
3:C:87:GLU:O	3:C:88:GLU:HB2	1.94	0.66
1:A:2:ASP:HA	1:A:5:GLU:HG2	1.78	0.65
1:A:212:PRO:O	1:A:213:ASN:CB	2.45	0.65
1:A:212:PRO:O	1:A:213:ASN:HB3	1.96	0.65
2:B:93:PRO:HG3	2:B:269:THR:C	2.22	0.64
3:C:112:GLU:HG3	3:C:113:PRO:HD2	1.78	0.64
2:B:97:LEU:HD13	2:B:97:LEU:O	1.98	0.64
3:C:17:ARG:HG2	3:C:19:GLY:H	1.63	0.64
1:A:7:GLU:HB2	1:A:10:ASN:HB3	1.81	0.63
2:B:68:GLU:HG3	2:B:71:HIS:CD2	2.33	0.63
3:C:53:ARG:HD2	3:C:53:ARG:N	2.13	0.63
3:C:104:GLU:O	3:C:106:LYS:N	2.32	0.63
3:C:104:GLU:C	3:C:106:LYS:H	2.07	0.63
2:B:18:LEU:HD22	3:C:75:VAL:HG13	1.81	0.62
1:A:14:CYS:HB3	1:A:96:VAL:HG13	1.81	0.62
1:A:119:HIS:ND1	1:A:150:TYR:CE1	2.67	0.62
3:C:75:VAL:HG23	3:C:97:LEU:O	1.99	0.62
3:C:105:ASP:C	3:C:106:LYS:HD2	2.24	0.62
3:C:89:LYS:O	3:C:91:GLU:N	2.33	0.62
2:B:160:THR:O	2:B:163:TRP:HB3	2.00	0.62
3:C:28:LYS:HE2	3:C:30:HIS:NE2	2.15	0.62
3:C:112:GLU:O	3:C:114:LEU:N	2.32	0.62
1:A:212:PRO:O	1:A:213:ASN:ND2	2.33	0.61
2:B:63:ILE:HA	2:B:87:PHE:O	2.00	0.61
2:B:271:LYS:C	2:B:273:ASN:H	2.08	0.61
3:C:49:THR:HB	3:C:50:PRO:HD3	1.83	0.60
1:A:126:ILE:HD13	1:A:139:VAL:HG11	1.82	0.60
2:B:164:LEU:O	2:B:168:VAL:HG22	2.01	0.60
3:C:58:GLY:C	3:C:59:LEU:HG	2.27	0.59
1:A:30:VAL:HG21	1:A:57:LEU:HD21	1.83	0.59
2:B:60:LEU:HD23	2:B:91:MSE:HG2	1.85	0.59
1:A:10:ASN:HB2	3:C:53:ARG:NH1	2.17	0.58
1:A:111:LYS:O	4:A:339:HOH:O	2.17	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:222:ARG:HG3	1:A:249:GLU:OE2	2.02	0.58
2:B:91:MSE:O	2:B:92:ASP:CB	2.51	0.58
3:C:27:ALA:HB2	3:C:94:ILE:CB	2.34	0.58
3:C:28:LYS:HE2	3:C:30:HIS:CE1	2.39	0.57
1:A:85:GLU:HG3	1:A:86:ASP:N	2.19	0.57
1:A:1:MSE:O	1:A:3:LEU:N	2.37	0.57
3:C:75:VAL:HG13	3:C:76:PRO:HD2	1.86	0.57
2:B:50:THR:HB	3:C:142:VAL:HG11	1.85	0.57
2:B:10:LEU:O	2:B:11:ALA:HB2	2.04	0.57
2:B:101:LEU:CD2	2:B:178:ASN:ND2	2.68	0.56
2:B:101:LEU:CD2	2:B:178:ASN:HD22	2.19	0.56
2:B:98:LEU:HD12	2:B:101:LEU:HD23	1.86	0.56
3:C:136:LEU:C	3:C:138:GLY:H	2.14	0.56
1:A:65:VAL:HG13	1:A:83:MSE:HE2	1.88	0.55
3:C:18:PRO:HA	3:C:21:LEU:HD21	1.88	0.55
3:C:112:GLU:HG3	3:C:113:PRO:CD	2.36	0.55
3:C:21:LEU:H	3:C:21:LEU:CD2	2.17	0.55
3:C:66:ARG:CZ	3:C:104:GLU:OE2	2.55	0.55
3:C:107:ALA:O	3:C:108:ASP:OD1	2.25	0.55
3:C:110:ALA:O	3:C:111:GLN:HG3	2.07	0.55
1:A:42:LEU:HD12	1:A:210:GLY:HA2	1.89	0.54
1:A:27:GLU:OE2	1:A:55:SER:HB3	2.08	0.54
1:A:122:ALA:O	1:A:126:ILE:HG13	2.08	0.53
2:B:60:LEU:HB3	2:B:91:MSE:HB3	1.90	0.53
2:B:68:GLU:HG3	2:B:71:HIS:HD2	1.73	0.53
2:B:17:PHE:CD2	3:C:84:MSE:HG3	2.42	0.53
1:A:1:MSE:C	1:A:3:LEU:H	2.17	0.52
2:B:10:LEU:O	2:B:11:ALA:CB	2.56	0.52
1:A:7:GLU:C	1:A:9:ASP:N	2.67	0.52
3:C:86:THR:HG23	3:C:87:GLU:HG3	1.91	0.52
1:A:14:CYS:HB3	1:A:96:VAL:CG1	2.40	0.52
2:B:66:PHE:HB3	2:B:85:LEU:HB3	1.92	0.52
3:C:138:GLY:O	3:C:139:LEU:HG	2.09	0.52
1:A:231:PHE:HB3	1:A:235:VAL:HG21	1.91	0.52
2:B:170:GLN:O	2:B:173:VAL:HG12	2.09	0.52
1:A:71:LEU:HD23	1:A:71:LEU:H	1.74	0.51
2:B:52:LEU:HB3	2:B:61:PHE:HB2	1.93	0.51
2:B:16:VAL:HG11	2:B:73:TRP:CE2	2.45	0.51
1:A:197:VAL:O	1:A:201:GLN:HG3	2.11	0.50
2:B:169:ASN:HA	2:B:172:VAL:HB	1.92	0.50
3:C:85:VAL:HG13	3:C:85:VAL:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:59:ASP:O	1:A:63:LEU:HB2	2.12	0.50
2:B:163:TRP:CE3	2:B:267:LEU:HD22	2.47	0.50
1:A:119:HIS:CE1	1:A:150:TYR:CE1	3.00	0.50
3:C:18:PRO:HA	3:C:21:LEU:CD2	2.41	0.50
3:C:103:ALA:CB	3:C:108:ASP:HB3	2.41	0.50
2:B:162:LYS:HG2	2:B:165:GLU:OE1	2.12	0.50
3:C:141:THR:HG22	3:C:142:VAL:H	1.75	0.50
1:A:7:GLU:C	1:A:9:ASP:H	2.18	0.49
1:A:86:ASP:C	1:A:88:ASP:H	2.20	0.49
1:A:231:PHE:CE1	1:A:246:LEU:HG	2.47	0.49
1:A:59:ASP:O	1:A:63:LEU:CB	2.61	0.49
1:A:203:LEU:HD12	1:A:204:ASP:H	1.77	0.49
1:A:27:GLU:O	1:A:135:ASN:HB3	2.13	0.48
1:A:251:GLU:HA	1:A:251:GLU:OE1	2.12	0.48
1:A:222:ARG:NH1	1:A:222:ARG:CG	2.61	0.48
1:A:158:PHE:N	1:A:159:PRO:HD3	2.28	0.48
2:B:67:LYS:HB3	2:B:84:LEU:CD2	2.43	0.48
3:C:53:ARG:CG	3:C:53:ARG:NH1	2.68	0.48
3:C:142:VAL:N	3:C:143:PRO:HD2	2.29	0.48
1:A:143:THR:HG22	1:A:167:ALA:HA	1.96	0.48
1:A:1:MSE:N	4:A:340:HOH:O	2.47	0.48
3:C:16:LEU:HD23	3:C:16:LEU:O	2.13	0.48
1:A:149:THR:O	1:A:153:ARG:HG2	2.14	0.47
2:B:273:ASN:HA	2:B:276:MSE:HB3	1.96	0.47
2:B:178:ASN:OD1	2:B:178:ASN:N	2.45	0.47
3:C:20:SER:OG	3:C:92:GLY:N	2.48	0.47
1:A:103:SER:O	1:A:107:LEU:HB2	2.15	0.47
2:B:55:MSE:C	2:B:57:LEU:H	2.22	0.47
1:A:59:ASP:C	1:A:63:LEU:HB2	2.40	0.47
2:B:15:LEU:HB2	3:C:84:MSE:HB2	1.95	0.47
3:C:105:ASP:O	3:C:106:LYS:HD2	2.13	0.47
3:C:114:LEU:HD13	3:C:114:LEU:HA	1.55	0.46
2:B:45:SER:O	2:B:46:GLY:C	2.58	0.46
2:B:21:GLU:HG3	3:C:78:GLY:C	2.40	0.46
1:A:63:LEU:HB3	1:A:64:LYS:H	1.66	0.46
1:A:252:ASP:HB2	3:C:37:LEU:HD12	1.98	0.46
2:B:177:ALA:O	2:B:178:ASN:OD1	2.33	0.46
2:B:73:TRP:HB2	2:B:80:GLN:HB3	1.97	0.46
3:C:88:GLU:O	3:C:89:LYS:HG2	2.16	0.46
1:A:41:VAL:HB	1:A:239:TRP:HD1	1.79	0.46
3:C:66:ARG:NH1	3:C:104:GLU:OE2	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:88:GLU:C	3:C:89:LYS:HG2	2.40	0.46
1:A:220:TRP:O	1:A:224:HIS:HB2	2.16	0.45
1:A:24:CYS:SG	1:A:53:PRO:HD3	2.57	0.45
1:A:65:VAL:O	1:A:66:ALA:C	2.59	0.45
1:A:43:GLY:O	1:A:236:ARG:NH2	2.42	0.45
3:C:106:LYS:NZ	4:C:175:HOH:O	2.42	0.45
1:A:228:VAL:HG22	1:A:229:PHE:CD1	2.51	0.45
1:A:97:LEU:HD11	1:A:121:THR:HG21	1.98	0.45
3:C:106:LYS:O	3:C:107:ALA:C	2.59	0.45
1:A:140:PHE:HD1	1:A:164:THR:HB	1.82	0.44
2:B:148:ASN:ND2	2:B:276:MSE:SE	3.00	0.44
3:C:46:ARG:HB2	3:C:46:ARG:HH11	1.81	0.44
1:A:10:ASN:HB2	3:C:53:ARG:HH11	1.82	0.44
2:B:19:LEU:HD22	3:C:82:PHE:CD2	2.52	0.44
2:B:93:PRO:HB2	2:B:170:GLN:HE22	1.82	0.44
1:A:64:LYS:HB2	1:A:64:LYS:HE2	1.72	0.44
3:C:93:LEU:HD22	3:C:93:LEU:O	2.18	0.44
1:A:5:GLU:HG2	1:A:5:GLU:H	1.57	0.44
2:B:89:THR:HG21	2:B:100:TYR:CD1	2.52	0.44
1:A:2:ASP:O	1:A:3:LEU:C	2.60	0.44
3:C:142:VAL:N	3:C:143:PRO:CD	2.80	0.44
1:A:61:GLU:HG3	1:A:62:ALA:N	2.33	0.43
1:A:153:ARG:HG2	1:A:153:ARG:H	1.55	0.43
1:A:4:SER:O	1:A:7:GLU:HG2	2.18	0.43
1:A:119:HIS:O	1:A:120:ASP:C	2.61	0.43
1:A:86:ASP:HB2	1:A:89:PHE:HE2	1.84	0.43
3:C:139:LEU:C	3:C:139:LEU:HD12	2.43	0.43
1:A:13:ARG:HD3	1:A:95:ASP:OD2	2.18	0.43
2:B:16:VAL:HG22	2:B:85:LEU:CD1	2.49	0.43
2:B:68:GLU:CG	2:B:71:HIS:HD2	2.31	0.43
2:B:93:PRO:HG3	2:B:269:THR:O	2.19	0.43
2:B:65:VAL:O	3:C:139:LEU:HA	2.18	0.43
2:B:68:GLU:CG	2:B:71:HIS:CD2	3.00	0.43
1:A:39:GLY:HA3	1:A:211:TYR:HE2	1.84	0.43
3:C:61:ALA:O	3:C:68:LEU:HB2	2.18	0.43
1:A:99:PRO:HB2	1:A:233:GLN:O	2.18	0.43
1:A:40:PRO:HG2	1:A:187:ARG:NH2	2.33	0.43
1:A:51:TYR:O	1:A:52:CYS:HB3	2.18	0.43
2:B:93:PRO:HB2	2:B:170:GLN:HE21	1.80	0.43
3:C:89:LYS:NZ	3:C:89:LYS:HA	2.34	0.42
3:C:99:PHE:CD1	3:C:99:PHE:C	2.97	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:18:SER:O	1:A:19:PRO:C	2.62	0.42
3:C:34:CYS:SG	3:C:63:PHE:CE2	3.12	0.42
3:C:112:GLU:C	3:C:114:LEU:N	2.77	0.42
1:A:140:PHE:HA	1:A:164:THR:O	2.19	0.42
2:B:16:VAL:HG11	2:B:73:TRP:NE1	2.34	0.42
1:A:139:VAL:CG2	1:A:161:ILE:HD12	2.49	0.42
1:A:34:ASP:OD1	1:A:142:ASP:HB3	2.19	0.42
1:A:61:GLU:HB2	1:A:174:PRO:HG2	2.02	0.42
1:A:208:GLY:HA2	1:A:216:LYS:HD3	2.02	0.42
2:B:19:LEU:HD21	3:C:84:MSE:HE1	2.02	0.42
1:A:213:ASN:O	1:A:215:PRO:HD3	2.20	0.41
3:C:112:GLU:O	3:C:113:PRO:C	2.61	0.41
2:B:271:LYS:C	2:B:273:ASN:N	2.73	0.41
1:A:7:GLU:HB2	1:A:10:ASN:CB	2.49	0.41
3:C:17:ARG:NH1	3:C:20:SER:HB2	2.35	0.41
2:B:94:LEU:CB	2:B:170:GLN:HG2	2.49	0.41
3:C:112:GLU:OE1	3:C:112:GLU:HA	2.21	0.41
1:A:20:VAL:HA	1:A:21:PRO:HD3	1.96	0.41
1:A:41:VAL:O	1:A:41:VAL:HG22	2.21	0.41
1:A:255:TRP:CE3	3:C:34:CYS:HB3	2.55	0.41
2:B:19:LEU:HD12	2:B:88:ALA:HB3	2.02	0.41
1:A:189:LYS:HA	1:A:189:LYS:HD3	1.89	0.41
1:A:227:PRO:HG2	3:C:40:ARG:NH2	2.36	0.41
3:C:136:LEU:C	3:C:138:GLY:N	2.78	0.41
1:A:42:LEU:HD11	1:A:217:THR:HG23	2.02	0.41
1:A:57:LEU:HD23	1:A:57:LEU:HA	1.72	0.41
1:A:105:SER:HB3	1:A:112:TYR:HD2	1.86	0.41
2:B:16:VAL:HG22	2:B:16:VAL:O	2.20	0.41
2:B:65:VAL:HG12	2:B:86:HIS:CG	2.56	0.41
3:C:81:GLY:HA3	3:C:96:LYS:CE	2.51	0.41
3:C:81:GLY:HA3	3:C:96:LYS:HE2	2.03	0.41
1:A:18:SER:HB3	1:A:92:TRP:CE2	2.56	0.40
1:A:15:ARG:O	1:A:16:LEU:HD23	2.21	0.40
1:A:105:SER:HB3	1:A:112:TYR:CD2	2.56	0.40
1:A:47:TYR:HB3	1:A:125:LEU:HD12	2.04	0.40
1:A:108:GLY:N	4:A:324:HOH:O	2.47	0.40
1:A:166:LYS:HB2	1:A:169:ALA:HB2	2.04	0.40
1:A:203:LEU:CD1	1:A:204:ASP:H	2.34	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	252/301 (84%)	229 (91%)	20 (8%)	3 (1%)	10	34
2	B	132/332 (40%)	97 (74%)	21 (16%)	14 (11%)	0	1
3	C	109/166 (66%)	81 (74%)	16 (15%)	12 (11%)	0	1
All	All	493/799 (62%)	407 (83%)	57 (12%)	29 (6%)	1	4

All (29) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	11	ALA
2	B	46	GLY
2	B	92	ASP
2	B	156	SER
2	B	158	GLU
2	B	180	VAL
3	C	88	GLU
3	C	109	GLU
3	C	143	PRO
1	A	2	ASP
1	A	63	LEU
1	A	257	ASP
2	B	93	PRO
2	B	270	GLY
3	C	90	GLY
3	C	105	ASP
3	C	108	ASP
3	C	110	ALA
3	C	139	LEU
2	B	269	THR
2	B	47	GLU
2	B	150	LYS
2	B	272	LYS
3	C	94	ILE

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Mol	Chain	Res	Type
2	B	58	GLN
3	C	91	GLU
3	C	113	PRO
2	B	43	PRO
3	C	142	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	215/252 (85%)	175 (81%)	40 (19%)	1	5
2	B	82/286 (29%)	72 (88%)	10 (12%)	5	16
3	C	80/125 (64%)	54 (68%)	26 (32%)	0	1
All	All	377/663 (57%)	301 (80%)	76 (20%)	1	4

All (76) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	GLU
1	A	10	ASN
1	A	20	VAL
1	A	24	CYS
1	A	30	VAL
1	A	31	LEU
1	A	33	VAL
1	A	35	GLU
1	A	38	ARG
1	A	41	VAL
1	A	54	LEU
1	A	63	LEU
1	A	64	LYS
1	A	65	VAL
1	A	67	ASP
1	A	71	LEU
1	A	85	GLU

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Mol	Chain	Res	Type
1	A	94	LEU
1	A	95	ASP
1	A	107	LEU
1	A	114	LEU
1	A	138	GLN
1	A	148	GLU
1	A	149	THR
1	A	157	HIS
1	A	164	THR
1	A	168	LYS
1	A	172	LEU
1	A	192	LYS
1	A	200	LEU
1	A	201	GLN
1	A	203	LEU
1	A	213	ASN
1	A	222	ARG
1	A	224	HIS
1	A	228	VAL
1	A	236	ARG
1	A	241	THR
1	A	253	VAL
1	A	257	ASP
2	B	16	VAL
2	B	19	LEU
2	B	36	LEU
2	B	57	LEU
2	B	63	ILE
2	B	67	LYS
2	B	97	LEU
2	B	172	VAL
2	B	178	ASN
2	B	274	SER
3	C	14	ILE
3	C	16	LEU
3	C	17	ARG
3	C	21	LEU
3	C	28	LYS
3	C	46	ARG
3	C	53	ARG
3	C	57	ASP
3	C	59	LEU

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Mol	Chain	Res	Type
3	C	60	GLN
3	C	62	SER
3	C	73	VAL
3	C	84	MSE
3	C	88	GLU
3	C	89	LYS
3	C	93	LEU
3	C	96	LYS
3	C	104	GLU
3	C	105	ASP
3	C	106	LYS
3	C	108	ASP
3	C	112	GLU
3	C	114	LEU
3	C	115	GLU
3	C	139	LEU
3	C	141	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	113	ASN
1	A	135	ASN
1	A	201	GLN
1	A	224	HIS
2	B	14	GLN
2	B	58	GLN
2	B	70	HIS
2	B	148	ASN
2	B	170	GLN
3	C	98	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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	251/301 (83%)	0.24	19 (7%)	20 16	19, 36, 114, 151	0
2	B	137/332 (41%)	1.30	34 (24%)	2 1	30, 106, 149, 159	0
3	C	112/166 (67%)	0.69	19 (16%)	4 3	19, 47, 130, 155	0
All	All	500/799 (62%)	0.63	72 (14%)	6 5	19, 46, 136, 159	0

All (72) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	87	GLU	6.6
3	C	144	GLY	5.2
1	A	65	VAL	5.0
3	C	137	TRP	5.0
1	A	3	LEU	4.8
1	A	62	ALA	4.8
2	B	146	GLU	4.5
2	B	45	SER	4.2
3	C	141	THR	4.0
1	A	6	LEU	3.9
2	B	102	LEU	3.8
1	A	8	ARG	3.8
1	A	66	ALA	3.7
2	B	271	LYS	3.6
1	A	214	ASP	3.6
2	B	38	VAL	3.5
3	C	90	GLY	3.5
3	C	139	LEU	3.4
2	B	279	ALA	3.4
2	B	161	LEU	3.4
2	B	33	SER	3.4
2	B	158	GLU	3.3
1	A	63	LEU	3.3

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Mol	Chain	Res	Type	RSRZ
2	B	31	LYS	3.3
2	B	147	VAL	3.2
2	B	160	THR	3.2
3	C	138	GLY	3.2
2	B	282	ALA	3.1
2	B	36	LEU	3.0
2	B	181	ASN	2.9
3	C	86	THR	2.9
3	C	88	GLU	2.8
3	C	136	LEU	2.8
1	A	258	SER	2.7
3	C	111	GLN	2.7
2	B	37	PHE	2.6
1	A	71	LEU	2.6
2	B	57	LEU	2.6
2	B	278	ALA	2.6
2	B	43	PRO	2.6
3	C	21	LEU	2.6
2	B	156	SER	2.5
3	C	143	PRO	2.5
2	B	270	GLY	2.5
1	A	175	VAL	2.4
2	B	174	ALA	2.4
2	B	267	LEU	2.4
3	C	46	ARG	2.4
3	C	140	GLU	2.4
2	B	157	SER	2.4
1	A	10	ASN	2.3
2	B	11	ALA	2.3
2	B	272	LYS	2.3
3	C	58	GLY	2.3
1	A	9	ASP	2.3
1	A	2	ASP	2.3
1	A	256	GLU	2.2
3	C	13	ARG	2.2
2	B	100	TYR	2.2
2	B	274	SER	2.2
2	B	97	LEU	2.2
1	A	74	ASN	2.2
2	B	39	LYS	2.2
2	B	101	LEU	2.2
3	C	15	HIS	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	88	ASP	2.1
1	A	111	LYS	2.0
2	B	99	HIS	2.0
2	B	95	PHE	2.0
2	B	155	TYR	2.0
1	A	64	LYS	2.0
3	C	109	GLU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.