



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 06:53 PM UTC

PDB ID : 3G65 / pdb_00003g65
Title : Crystal Structure of the Human Rad9-Rad1-Hus1 DNA Damage Checkpoint Complex
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Deposited on : 2009-02-06
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
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

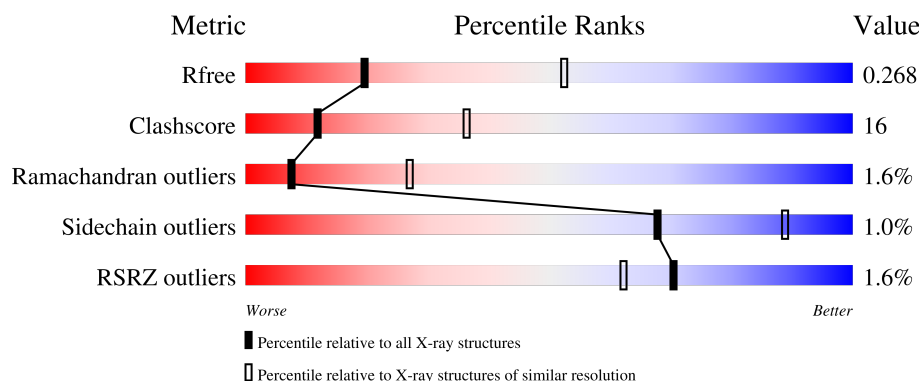
1 Overall quality at a glance

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	296	% 58% 27% • 14%
2	B	282	% 55% 36% • 7%
3	C	280	% 60% 30% • 8%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6150 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 Cell cycle checkpoint control protein RAD9A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	254	1958	1247	340	355	16	0	0	0

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	271	GLY	-	expression tag	UNP Q99638
A	272	THR	-	expression tag	UNP Q99638
A	273	THR	-	expression tag	UNP Q99638
A	274	SER	-	expression tag	UNP Q99638
A	275	THR	-	expression tag	UNP Q99638
A	276	SER	-	expression tag	UNP Q99638
A	277	LEU	-	expression tag	UNP Q99638
A	278	GLU	-	expression tag	UNP Q99638
A	279	VAL	-	expression tag	UNP Q99638
A	280	LEU	-	expression tag	UNP Q99638
A	281	PHE	-	expression tag	UNP Q99638
A	282	GLN	-	expression tag	UNP Q99638
A	283	GLY	-	expression tag	UNP Q99638
A	284	PRO	-	expression tag	UNP Q99638
A	285	LEU	-	expression tag	UNP Q99638
A	286	SER	-	expression tag	UNP Q99638
A	287	GLY	-	expression tag	UNP Q99638
A	288	SER	-	expression tag	UNP Q99638
A	289	GLY	-	expression tag	UNP Q99638
A	290	GLY	-	expression tag	UNP Q99638
A	291	HIS	-	expression tag	UNP Q99638
A	292	HIS	-	expression tag	UNP Q99638
A	293	HIS	-	expression tag	UNP Q99638
A	294	HIS	-	expression tag	UNP Q99638
A	295	HIS	-	expression tag	UNP Q99638
A	296	HIS	-	expression tag	UNP Q99638

- Molecule 2 is a protein called Cell cycle checkpoint protein RAD1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	261	Total	C	N	O	S	0	0	0
			2058	1306	338	396	18			

- Molecule 3 is a protein called Checkpoint protein HUS1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	258	Total	C	N	O	S	0	0	0
			2052	1312	348	376	16			

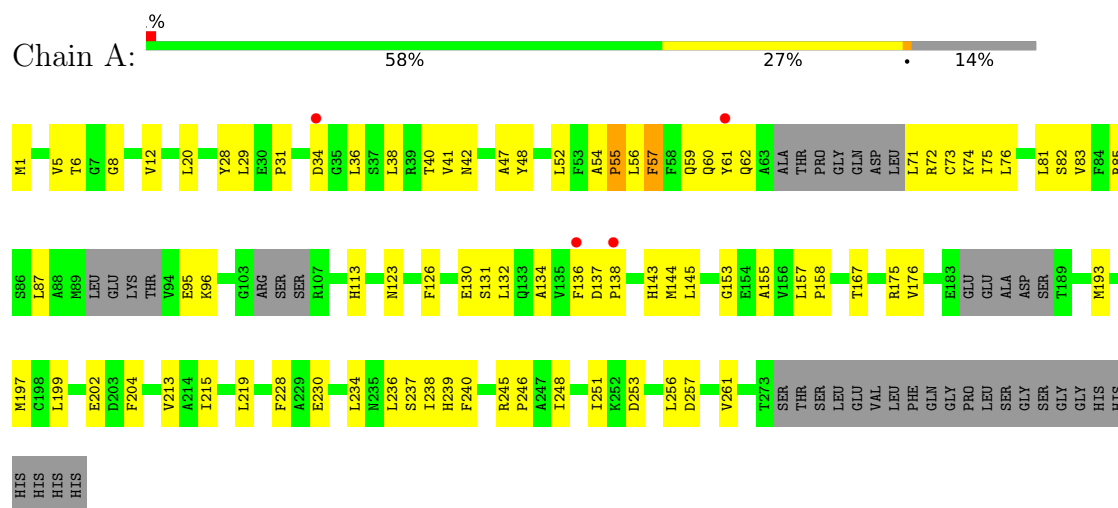
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	33	Total	O	0	0
			33	33		
4	B	24	Total	O	0	0
			24	24		
4	C	25	Total	O	0	0
			25	25		

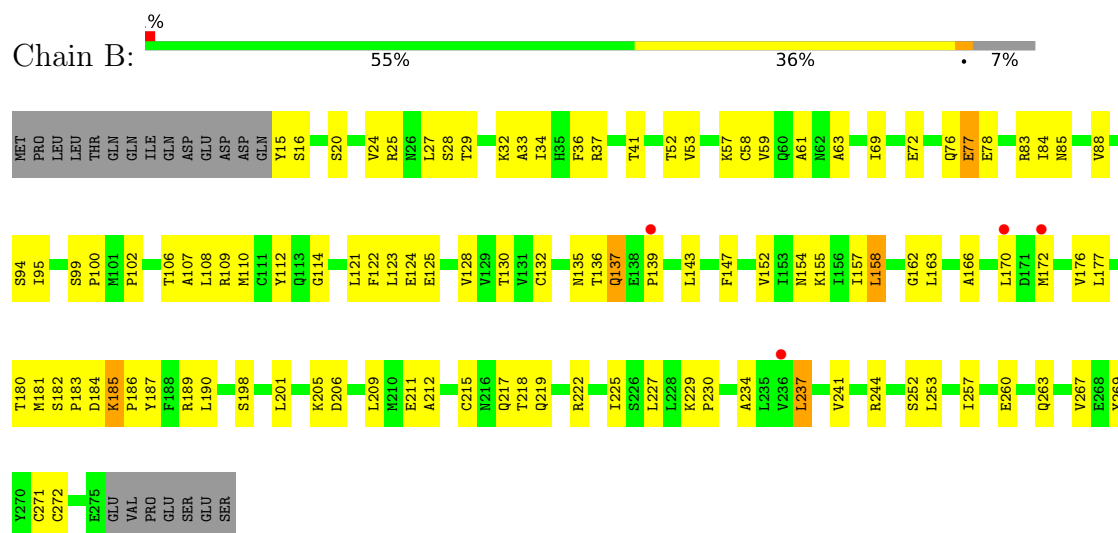
3 Residue-property plots [i](#)

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: Cell cycle checkpoint control protein RAD9A

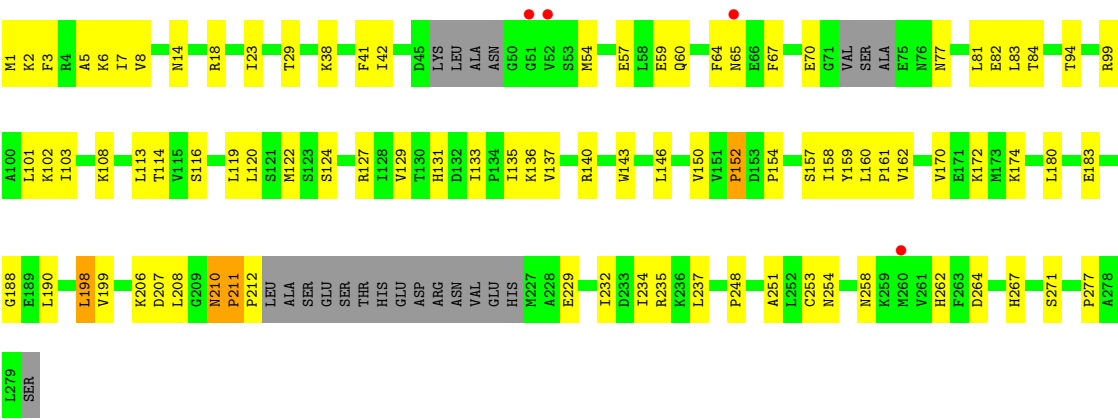


- Molecule 2: Cell cycle checkpoint protein RAD1



- Molecule 3: Checkpoint protein HUS1





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	73.25Å 70.67Å 83.15Å 90.00° 99.51° 90.00°	Depositor
Resolution (Å)	82.00 – 2.90 82.00 – 2.90	Depositor EDS
% Data completeness (in resolution range)	96.5 (82.00-2.90) 97.4 (82.00-2.90)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.58 (at 2.86Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.221 , 0.272 0.218 , 0.268	Depositor DCC
R_{free} test set	942 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	76.7	Xtriage
Anisotropy	0.202	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 87.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6150	wwPDB-VP
Average B, all atoms (Å ²)	87.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.28% 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	0.23	0/1990	0.66	0/2679
2	B	0.24	0/2096	0.68	1/2835 (0.0%)
3	C	0.24	0/2086	0.66	0/2818
All	All	0.24	0/6172	0.67	1/8332 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	77	GLU	CB-CA-C	-5.15	110.23	117.23

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1958	0	1988	59	0
2	B	2058	0	2038	80	0
3	C	2052	0	2116	63	0
4	A	33	0	0	1	0
4	B	24	0	0	1	0
4	C	25	0	0	0	0
All	All	6150	0	6142	193	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 16.

All (193) 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)
3:C:211:PRO:HB2	3:C:212:PRO:HD3	1.39	1.04
2:B:185:LYS:HB2	2:B:186:PRO:HD3	1.40	1.04
2:B:59:VAL:HG21	2:B:227:LEU:HB3	1.54	0.89
2:B:20:SER:HB3	2:B:72:GLU:HB3	1.59	0.85
1:A:42:ASN:HB2	1:A:47:ALA:HB3	1.58	0.83
3:C:57:GLU:HB3	3:C:271:SER:HB3	1.60	0.82
3:C:232:ILE:HD11	3:C:237:LEU:HD12	1.64	0.79
1:A:96:LYS:HB2	1:A:113:HIS:HB2	1.66	0.78
2:B:106:THR:HG22	2:B:125:GLU:HB2	1.70	0.73
2:B:147:PHE:HE1	2:B:244:ARG:HD2	1.53	0.72
2:B:57:LYS:HD2	2:B:143:LEU:HD13	1.72	0.71
3:C:38:LYS:HE2	3:C:59:GLU:HG2	1.73	0.70
2:B:24:VAL:HG22	2:B:108:LEU:HB2	1.74	0.70
2:B:162:GLY:HA3	2:B:209:LEU:HD21	1.74	0.69
1:A:253:ASP:HB3	1:A:256:LEU:HB3	1.74	0.68
2:B:135:ASN:HB2	3:C:198:LEU:HD12	1.76	0.68
3:C:174:LYS:HE3	3:C:235:ARG:HG2	1.78	0.65
2:B:99:SER:N	2:B:100:PRO:HD3	2.12	0.65
1:A:158:PRO:HG2	1:A:197:MET:HE2	1.80	0.64
1:A:41:VAL:HG23	1:A:132:LEU:HD13	1.79	0.64
3:C:183:GLU:HG2	3:C:229:GLU:HG2	1.78	0.64
3:C:159:TYR:HE1	3:C:211:PRO:HD2	1.60	0.64
1:A:83:VAL:HA	1:A:87:LEU:HD12	1.80	0.63
2:B:34:ILE:HG13	2:B:63:ALA:HB2	1.80	0.63
1:A:75:ILE:HG12	1:A:126:PHE:HB3	1.81	0.62
2:B:95:ILE:HA	3:C:172:LYS:HG3	1.81	0.62
2:B:176:VAL:HG21	2:B:222:ARG:HB3	1.82	0.62
2:B:185:LYS:HB2	2:B:186:PRO:CD	2.23	0.62
3:C:1:MET:HE1	3:C:137:VAL:HG21	1.80	0.62
1:A:61:TYR:HB2	1:A:71:LEU:HD12	1.82	0.61
3:C:8:VAL:HG13	3:C:99:ARG:HH22	1.65	0.61
1:A:155:ALA:HB2	1:A:199:LEU:HD11	1.83	0.61
2:B:24:VAL:HG23	2:B:106:THR:O	2.00	0.61
2:B:185:LYS:CB	2:B:186:PRO:HD3	2.24	0.60
3:C:29:THR:HG22	3:C:84:THR:HG22	1.82	0.60
2:B:225:ILE:O	2:B:229:LYS:HG3	2.02	0.60
2:B:217:GLN:HE21	2:B:219:GLN:NE2	2.00	0.59
3:C:99:ARG:HB2	3:C:120:LEU:HG	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:147:PHE:CE1	2:B:244:ARG:HD2	2.37	0.57
2:B:83:ARG:HB3	2:B:137:GLN:O	2.04	0.57
2:B:253:LEU:HB2	2:B:269:TYR:HB2	1.87	0.57
3:C:103:ILE:HD11	3:C:113:LEU:HD21	1.87	0.57
1:A:136:PHE:CE1	1:A:261:VAL:HG21	2.39	0.57
2:B:59:VAL:CG2	2:B:227:LEU:HB3	2.32	0.57
2:B:83:ARG:HD3	2:B:139:PRO:HD2	1.87	0.56
3:C:160:LEU:HB3	3:C:251:ALA:HB1	1.86	0.56
1:A:52:LEU:HB3	1:A:257:ASP:HB3	1.88	0.56
1:A:143:HIS:CE1	1:A:213:VAL:HG21	2.41	0.56
2:B:88:VAL:HG11	3:C:199:VAL:HG21	1.87	0.56
1:A:6:THR:HG22	1:A:96:LYS:HG2	1.88	0.55
3:C:140:ARG:HA	3:C:143:TRP:CD1	2.41	0.55
3:C:158:ILE:HA	3:C:210:ASN:HB3	1.89	0.55
3:C:38:LYS:HG2	3:C:59:GLU:HA	1.87	0.55
2:B:112:TYR:CE1	2:B:136:THR:HG21	2.41	0.54
2:B:217:GLN:HE21	2:B:219:GLN:HE21	1.52	0.54
2:B:15:TYR:O	2:B:16:SER:HB3	2.08	0.54
3:C:5:ALA:HB3	3:C:103:ILE:HG23	1.90	0.54
2:B:77:GLU:HG2	2:B:78:GLU:HG3	1.89	0.54
2:B:183:PRO:HG3	2:B:215:CYS:O	2.09	0.53
2:B:109:ARG:HB2	2:B:122:PHE:HB2	1.89	0.53
3:C:23:ILE:HD12	3:C:41:PHE:HB3	1.90	0.53
2:B:229:LYS:HB2	2:B:230:PRO:HD3	1.89	0.53
1:A:153:GLY:HA3	1:A:230:GLU:OE2	2.08	0.53
1:A:130:GLU:HG2	1:A:131:SER:H	1.74	0.53
3:C:211:PRO:HB2	3:C:212:PRO:CD	2.18	0.53
2:B:28:SER:O	2:B:32:LYS:HG3	2.09	0.52
3:C:157:SER:O	3:C:210:ASN:HB2	2.08	0.52
2:B:37:ARG:HB2	2:B:53:VAL:CG1	2.39	0.52
2:B:166:ALA:O	2:B:170:LEU:HB2	2.10	0.51
3:C:5:ALA:HB1	3:C:64:PHE:CD1	2.46	0.51
2:B:34:ILE:HG23	2:B:61:ALA:HB3	1.92	0.51
3:C:262:HIS:CE1	3:C:264:ASP:HB2	2.46	0.51
1:A:62:GLN:H	1:A:71:LEU:HD11	1.75	0.51
1:A:138:PRO:HG3	1:A:248:ILE:HG21	1.92	0.51
1:A:234:LEU:HD12	1:A:251:ILE:HD11	1.93	0.50
3:C:102:LYS:HB2	3:C:116:SER:OG	2.12	0.50
1:A:134:ALA:HB3	1:A:136:PHE:CZ	2.47	0.50
2:B:152:VAL:HG21	2:B:244:ARG:HD3	1.93	0.50
1:A:158:PRO:HA	3:C:94:THR:HG21	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:75:ILE:HG22	1:A:76:LEU:N	2.26	0.50
2:B:170:LEU:O	2:B:172:MET:HG2	2.12	0.49
1:A:59:GLN:HG2	1:A:60:GLN:N	2.26	0.49
2:B:184:ASP:O	2:B:185:LYS:C	2.56	0.49
2:B:84:ILE:HG12	2:B:85:ASN:N	2.27	0.49
1:A:54:ALA:O	1:A:56:LEU:N	2.45	0.49
3:C:152:PRO:O	3:C:154:PRO:HD3	2.13	0.49
3:C:82:GLU:HB3	3:C:136:LYS:HB3	1.95	0.48
2:B:36:PHE:HB3	2:B:230:PRO:HB3	1.95	0.48
2:B:225:ILE:HG23	2:B:229:LYS:HE3	1.96	0.48
2:B:107:ALA:O	2:B:123:LEU:HA	2.12	0.48
2:B:180:THR:HG23	2:B:189:ARG:HB3	1.96	0.48
1:A:28:TYR:CE1	1:A:74:LYS:HD3	2.48	0.48
2:B:244:ARG:HB3	2:B:252:SER:HB3	1.95	0.48
1:A:215:ILE:HD13	1:A:240:PHE:HB2	1.97	0.47
2:B:121:LEU:HB2	2:B:132:CYS:HB2	1.97	0.47
2:B:182:SER:O	2:B:187:TYR:HA	2.14	0.47
3:C:154:PRO:HB3	3:C:254:ASN:HB3	1.96	0.47
2:B:25:ARG:HD3	2:B:260:GLU:OE2	2.15	0.47
2:B:108:LEU:HD13	2:B:123:LEU:HD21	1.96	0.46
3:C:7:ILE:HD12	3:C:101:LEU:HD23	1.96	0.46
3:C:14:ASN:O	3:C:18:ARG:HG3	2.15	0.46
1:A:81:LEU:O	1:A:85:ARG:HG3	2.15	0.46
1:A:202:GLU:OE1	3:C:124:SER:HB3	2.15	0.46
2:B:29:THR:HG22	2:B:237:LEU:HD21	1.98	0.46
2:B:41:THR:HB	2:B:52:THR:HG23	1.97	0.46
2:B:58:CYS:O	2:B:271:CYS:HA	2.16	0.46
3:C:237:LEU:HD23	3:C:237:LEU:O	2.15	0.46
2:B:58:CYS:SG	2:B:272:CYS:HB2	2.55	0.46
1:A:158:PRO:HB3	3:C:129:VAL:HG11	1.98	0.46
1:A:5:VAL:HG22	1:A:57:PHE:CD2	2.51	0.46
1:A:20:LEU:HD22	1:A:40:THR:HG22	1.98	0.46
1:A:176:VAL:HG21	1:A:204:PHE:CZ	2.51	0.46
1:A:228:PHE:CZ	1:A:256:LEU:HD21	2.51	0.46
2:B:94:SER:O	3:C:172:LYS:HE3	2.15	0.46
3:C:81:LEU:HD13	3:C:135:ILE:HG12	1.98	0.45
1:A:95:GLU:HG2	1:A:113:HIS:O	2.16	0.45
3:C:8:VAL:HG13	3:C:99:ARG:NH2	2.30	0.45
2:B:34:ILE:HD11	2:B:267:VAL:HG13	1.97	0.45
3:C:83:LEU:HB3	3:C:135:ILE:HD12	1.99	0.45
3:C:211:PRO:CB	3:C:212:PRO:HD3	2.26	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:57:LYS:HB3	2:B:143:LEU:HD22	1.98	0.45
2:B:69:ILE:HA	2:B:263:GLN:OE1	2.16	0.45
2:B:170:LEU:HD21	2:B:177:LEU:HD22	1.98	0.44
3:C:114:THR:HA	3:C:131:HIS:O	2.17	0.44
1:A:1:MET:SD	1:A:31:PRO:HG3	2.58	0.44
1:A:36:LEU:HG	1:A:38:LEU:CD1	2.48	0.44
1:A:136:PHE:O	1:A:137:ASP:HB2	2.18	0.44
1:A:145:LEU:HD11	1:A:238:ILE:HD12	1.99	0.44
2:B:157:ILE:O	2:B:211:GLU:HB3	2.16	0.44
3:C:6:LYS:HD2	3:C:102:LYS:HE2	2.00	0.44
2:B:237:LEU:O	2:B:237:LEU:HD23	2.17	0.44
3:C:160:LEU:HD13	3:C:253:CYS:HB2	2.00	0.44
2:B:24:VAL:HG21	2:B:123:LEU:HD22	1.99	0.44
3:C:119:LEU:HB2	3:C:127:ARG:HB2	2.00	0.44
2:B:170:LEU:HG	2:B:225:ILE:HD11	1.99	0.44
2:B:218:THR:HG22	4:B:345:HOH:O	2.18	0.44
1:A:157:LEU:HB2	1:A:158:PRO:HD3	1.99	0.43
3:C:2:LYS:HD2	3:C:70:GLU:CD	2.43	0.43
3:C:161:PRO:HG3	3:C:208:LEU:HD11	1.99	0.43
1:A:175:ARG:HB2	1:A:199:LEU:O	2.18	0.43
1:A:34:ASP:O	1:A:54:ALA:HA	2.18	0.43
1:A:48:TYR:CD2	1:A:136:PHE:HZ	2.37	0.43
3:C:188:GLY:HA3	3:C:208:LEU:O	2.19	0.43
3:C:180:LEU:C	3:C:180:LEU:HD23	2.43	0.43
1:A:8:GLY:O	1:A:12:VAL:HG23	2.18	0.43
1:A:144:MET:HE2	1:A:239:HIS:CD2	2.54	0.42
2:B:34:ILE:O	2:B:34:ILE:HG22	2.18	0.42
3:C:3:PHE:HE1	3:C:67:PHE:CZ	2.37	0.42
3:C:210:ASN:N	3:C:211:PRO:HD3	2.34	0.42
2:B:155:LYS:HE2	2:B:157:ILE:HD11	2.00	0.42
1:A:29:LEU:HB2	1:A:73:CYS:SG	2.59	0.42
3:C:23:ILE:HG13	3:C:54:MET:HE1	2.01	0.42
2:B:25:ARG:HH21	2:B:100:PRO:HD2	1.85	0.42
2:B:34:ILE:HG23	2:B:61:ALA:CB	2.50	0.42
2:B:95:ILE:HG21	2:B:130:THR:HG21	2.01	0.42
1:A:155:ALA:HA	1:A:197:MET:HE3	2.01	0.42
1:A:256:LEU:HD12	1:A:257:ASP:H	1.85	0.42
3:C:42:ILE:HD13	3:C:146:LEU:O	2.20	0.42
3:C:183:GLU:O	3:C:190:LEU:HD12	2.20	0.42
1:A:54:ALA:O	1:A:55:PRO:C	2.61	0.41
1:A:72:ARG:HD2	1:A:130:GLU:HA	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:95:GLU:HB2	4:A:377:HOH:O	2.19	0.41
3:C:122:MET:HA	3:C:122:MET:HE2	2.00	0.41
3:C:232:ILE:HG22	3:C:277:PRO:O	2.19	0.41
2:B:16:SER:HA	2:B:76:GLN:HB3	2.02	0.41
3:C:206:LYS:O	3:C:207:ASP:HB2	2.20	0.41
1:A:236:LEU:HD23	1:A:237:SER:N	2.35	0.41
1:A:176:VAL:HG22	1:A:199:LEU:HB2	2.02	0.41
2:B:157:ILE:HB	2:B:212:ALA:HB3	2.02	0.41
3:C:8:VAL:HB	3:C:65:ASN:ND2	2.36	0.41
2:B:124:GLU:HA	2:B:128:VAL:O	2.21	0.41
2:B:27:LEU:HD21	2:B:110:MET:HE1	2.02	0.41
3:C:248:PRO:HA	3:C:267:HIS:ND1	2.36	0.41
1:A:228:PHE:HZ	1:A:256:LEU:HD21	1.85	0.41
2:B:123:LEU:HB2	2:B:130:THR:CG2	2.51	0.41
2:B:190:LEU:HD12	2:B:201:LEU:HD23	2.02	0.41
1:A:193:MET:SD	3:C:133:ILE:HG12	2.61	0.41
2:B:205:LYS:O	2:B:206:ASP:HB2	2.21	0.41
2:B:237:LEU:HD22	2:B:257:ILE:HD13	2.03	0.41
1:A:123:ASN:HB2	2:B:198:SER:OG	2.21	0.40
2:B:84:ILE:HG12	2:B:85:ASN:H	1.86	0.40
3:C:6:LYS:HA	3:C:101:LEU:O	2.21	0.40
1:A:82:SER:HA	1:A:85:ARG:HD2	2.03	0.40
1:A:29:LEU:N	1:A:29:LEU:HD12	2.36	0.40
1:A:157:LEU:N	1:A:158:PRO:CD	2.83	0.40
1:A:245:ARG:HA	1:A:246:PRO:HD3	1.92	0.40
2:B:158:LEU:HD11	2:B:163:LEU:HD11	2.03	0.40
1:A:62:GLN:N	1:A:71:LEU:HD11	2.35	0.40
2:B:33:ALA:HB1	2:B:234:ALA:HB2	2.03	0.40
3:C:60:GLN:HG2	3:C:64:PHE:HD2	1.86	0.40
3:C:150:VAL:O	3:C:152:PRO:HD3	2.22	0.40
2:B:154:ASN:HB3	2:B:181:MET:HE2	2.03	0.40
3:C:170:VAL:HG13	3:C:234:ILE:CG1	2.51	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	244/296 (82%)	222 (91%)	20 (8%)	2 (1%)	16	44
2	B	259/282 (92%)	226 (87%)	30 (12%)	3 (1%)	10	34
3	C	250/280 (89%)	220 (88%)	23 (9%)	7 (3%)	4	15
All	All	753/858 (88%)	668 (89%)	73 (10%)	12 (2%)	7	27

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	55	PRO
2	B	185	LYS
3	C	152	PRO
3	C	258	ASN
3	C	108	LYS
1	A	57	PHE
3	C	77	ASN
3	C	198	LEU
3	C	211	PRO
2	B	114	GLY
3	C	210	ASN
2	B	102	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	215/250 (86%)	213 (99%)	2 (1%)	70	90
2	B	235/256 (92%)	231 (98%)	4 (2%)	53	82
3	C	238/257 (93%)	237 (100%)	1 (0%)	84	94
All	All	688/763 (90%)	681 (99%)	7 (1%)	68	89

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	167	THR
1	A	219	LEU
2	B	137	GLN
2	B	158	LEU
2	B	237	LEU
2	B	241	VAL
3	C	162	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	HIS
1	A	206	GLN
1	A	239	HIS
2	B	60	GLN
2	B	62	ASN
2	B	71	GLN
2	B	113	GLN
2	B	217	GLN
3	C	68	GLN
3	C	97	ASN
3	C	179	HIS
3	C	244	GLN
3	C	254	ASN
3	C	262	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	254/296 (85%)	0.09	4 (1%)	70 62	49, 86, 140, 184	0
2	B	261/282 (92%)	0.06	4 (1%)	72 64	43, 77, 135, 176	0
3	C	258/280 (92%)	0.08	4 (1%)	70 62	44, 80, 132, 159	0
All	All	773/858 (90%)	0.08	12 (1%)	70 62	43, 82, 135, 184	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	52	VAL	3.8
1	A	136	PHE	3.3
1	A	61	TYR	3.0
2	B	170	LEU	3.0
3	C	51	GLY	2.9
3	C	65	ASN	2.5
2	B	139	PRO	2.4
2	B	172	MET	2.2
2	B	236	VAL	2.1
3	C	260	MET	2.0
1	A	34	ASP	2.0
1	A	138	PRO	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.