



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 02:36 PM UTC

PDB ID : 1U4F / pdb_00001u4f
Title : Crystal Structure of Cytoplasmic Domains of IRK1 (Kir2.1) channel
Authors : Pegan, S.; Arrabit, C.; Zhou, W.; Kwiatkowski, W.; Slesinger, P.A.; Choe, S.
Deposited on : 2004-07-24
Resolution : 2.41 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

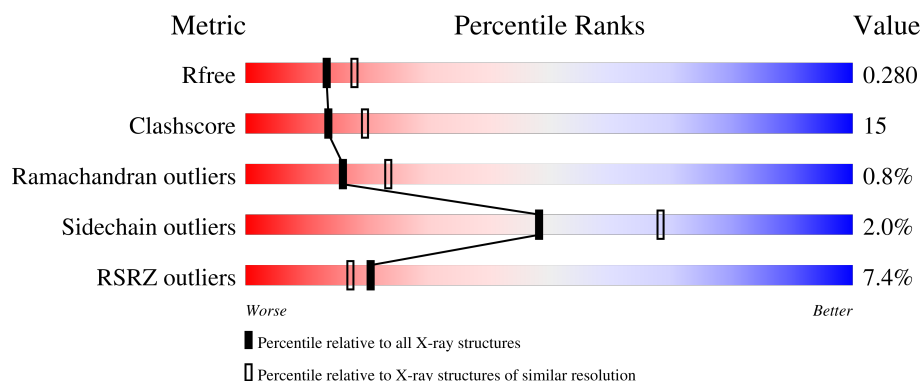
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.41 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	6062 (2.44-2.40)
Clashscore	190562	6562 (2.44-2.40)
Ramachandran outliers	187476	6481 (2.44-2.40)
Sidechain outliers	187428	6482 (2.44-2.40)
RSRZ outliers	180081	6066 (2.44-2.40)

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

Mol	Chain	Length	Quality of chain
1	A	270	 3% 51% 20% • 27%
1	B	270	 7% 54% 17% • 28%
1	C	270	 4% 48% 24% 28%
1	D	270	 6% 48% 24% • 27%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6533 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 Inward rectifier potassium channel 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	197	Total	C	N	O	S	0	0	0
			1590	1006	277	299	8			
1	B	195	Total	C	N	O	S	0	0	0
			1570	994	271	297	8			
1	C	195	Total	C	N	O	S	0	0	0
			1570	994	271	297	8			
1	D	197	Total	C	N	O	S	0	0	0
			1590	1012	273	297	8			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	35	PRO	-	cloning artifact	UNP P35561
A	36	ARG	-	cloning artifact	UNP P35561
A	37	GLY	-	cloning artifact	UNP P35561
A	38	SER	-	cloning artifact	UNP P35561
A	39	HIS	-	cloning artifact	UNP P35561
A	40	GLY	-	cloning artifact	UNP P35561
B	35	PRO	-	cloning artifact	UNP P35561
B	36	ARG	-	cloning artifact	UNP P35561
B	37	GLY	-	cloning artifact	UNP P35561
B	38	SER	-	cloning artifact	UNP P35561
B	39	HIS	-	cloning artifact	UNP P35561
B	40	GLY	-	cloning artifact	UNP P35561
C	35	PRO	-	cloning artifact	UNP P35561
C	36	ARG	-	cloning artifact	UNP P35561
C	37	GLY	-	cloning artifact	UNP P35561
C	38	SER	-	cloning artifact	UNP P35561
C	39	HIS	-	cloning artifact	UNP P35561
C	40	GLY	-	cloning artifact	UNP P35561
D	35	PRO	-	cloning artifact	UNP P35561
D	36	ARG	-	cloning artifact	UNP P35561
D	37	GLY	-	cloning artifact	UNP P35561

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Chain	Residue	Modelled	Actual	Comment	Reference
D	38	SER	-	cloning artifact	UNP P35561
D	39	HIS	-	cloning artifact	UNP P35561
D	40	GLY	-	cloning artifact	UNP P35561

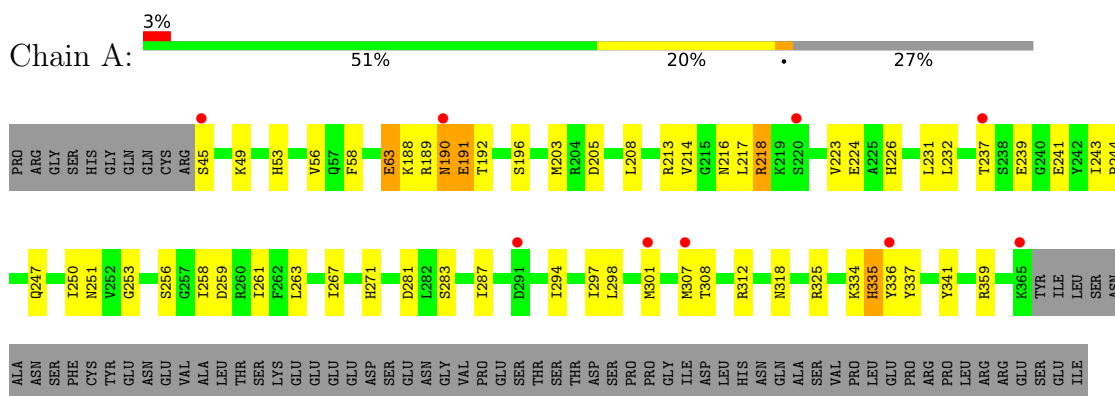
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	55	Total 55	O 55	0	0
2	B	48	Total 48	O 48	0	0
2	C	59	Total 59	O 59	0	0
2	D	51	Total 51	O 51	0	0

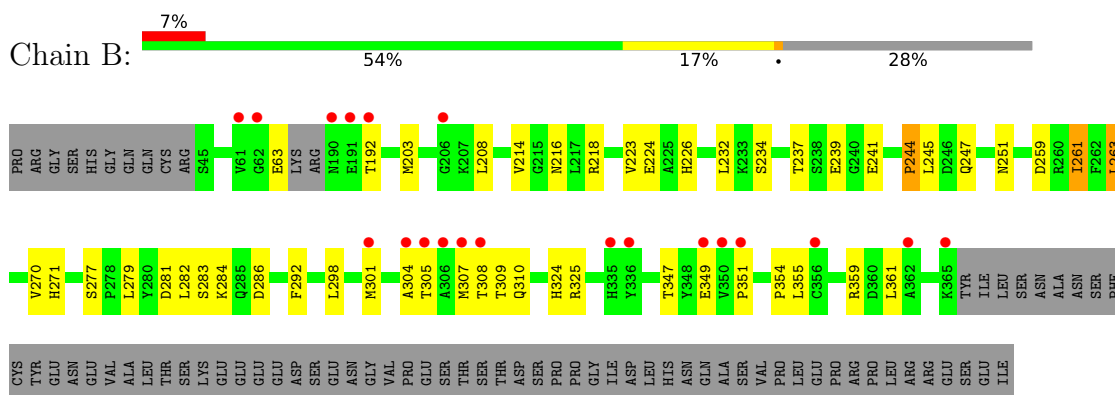
3 Residue-property plots

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.

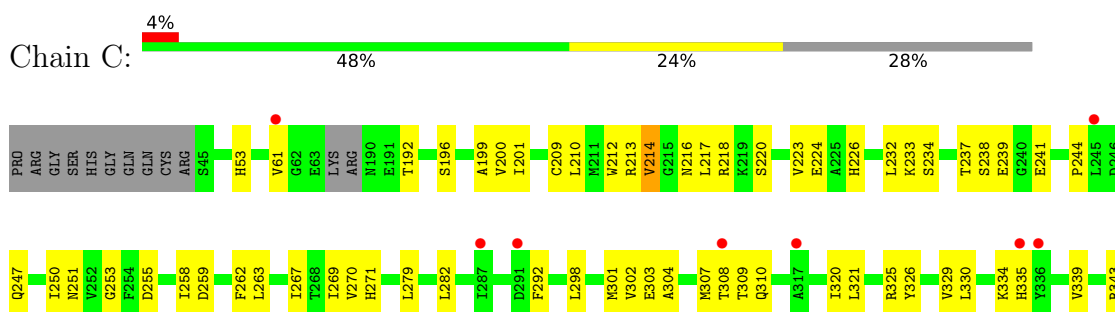
• Molecule 1: Inward rectifier potassium channel 2

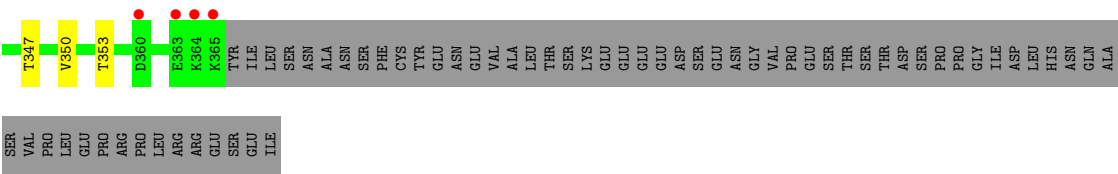


• Molecule 1: Inward rectifier potassium channel 2

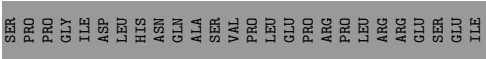
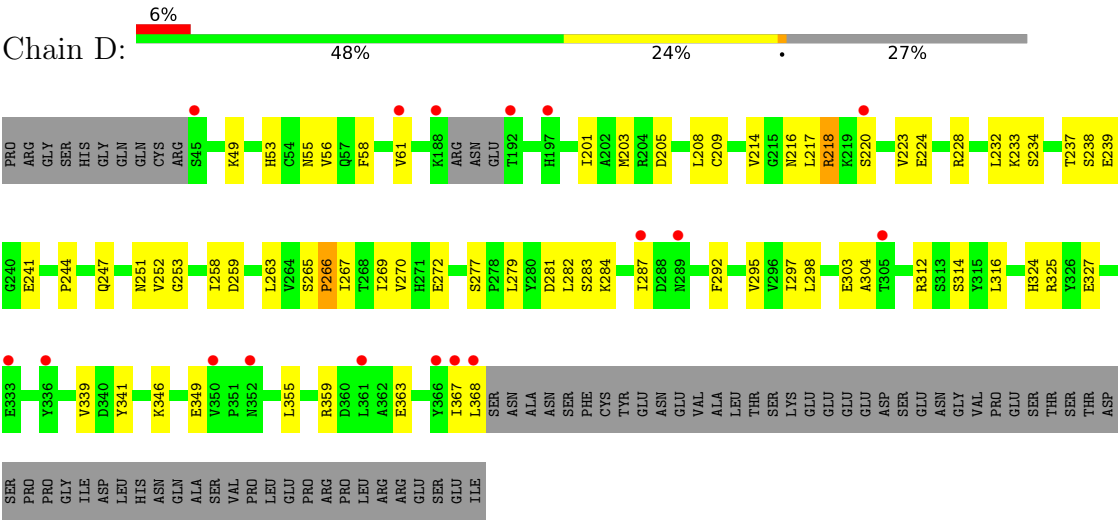


• Molecule 1: Inward rectifier potassium channel 2





● Molecule 1: Inward rectifier potassium channel 2



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	112.46Å 138.29Å 138.75Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.44 – 2.41 38.44 – 2.41	Depositor EDS
% Data completeness (in resolution range)	(Not available) (38.44-2.41) 92.1 (38.44-2.41)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.11 (at 2.39Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.225 , 0.281 0.226 , 0.280	Depositor DCC
R_{free} test set	1977 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	41.8	Xtriage
Anisotropy	0.849	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 40.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6533	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 12.48% 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	0.43	0/1622	0.89	4/2192 (0.2%)
1	B	0.44	0/1601	0.92	2/2164 (0.1%)
1	C	0.47	0/1601	0.92	6/2164 (0.3%)
1	D	0.43	0/1622	0.87	4/2192 (0.2%)
All	All	0.44	0/6446	0.90	16/8712 (0.2%)

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	259	ASP	N-CA-C	-8.42	103.00	113.18
1	B	259	ASP	N-CA-C	-7.50	104.11	113.18
1	A	259	ASP	N-CA-C	-7.28	104.37	113.18
1	C	259	ASP	N-CA-C	-7.14	104.38	113.23
1	B	270	VAL	N-CA-C	6.41	117.08	108.11
1	A	243	ILE	CA-C-N	5.77	127.06	119.84
1	A	243	ILE	C-N-CA	5.77	127.06	119.84
1	C	255	ASP	N-CA-C	5.66	118.21	111.71
1	C	270	VAL	N-CA-C	5.46	115.75	108.11
1	D	270	VAL	N-CA-C	5.38	115.64	108.11
1	D	251	ASN	N-CA-C	5.37	117.95	109.96
1	C	350	VAL	CA-C-N	5.34	125.67	119.47
1	C	350	VAL	C-N-CA	5.34	125.67	119.47
1	D	316	LEU	N-CA-C	-5.24	103.48	110.55
1	A	261	ILE	N-CA-C	5.12	116.97	108.99
1	C	353	THR	N-CA-C	5.03	116.85	109.71

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	1590	0	1570	51	0
1	B	1570	0	1543	46	0
1	C	1570	0	1543	50	0
1	D	1590	0	1575	54	0
2	A	55	0	0	0	0
2	B	48	0	0	1	0
2	C	59	0	0	2	0
2	D	51	0	0	2	0
All	All	6533	0	6231	183	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 15.

All (183) 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:B:203:MET:HE2	1:B:208:LEU:HD13	1.48	0.93
1:A:301:MET:HG2	1:A:308:THR:HG22	1.50	0.93
1:B:237:THR:HG23	1:B:239:GLU:H	1.38	0.87
1:A:188:LYS:HA	1:A:191:GLU:HG2	1.58	0.84
1:B:237:THR:HG22	1:B:241:GLU:H	1.43	0.83
1:A:223:VAL:HG12	1:A:224:GLU:HG3	1.62	0.82
1:B:301:MET:HG2	1:B:308:THR:HG22	1.63	0.81
1:D:297:ILE:HD12	1:D:312:ARG:HG2	1.64	0.80
1:B:237:THR:HG22	1:B:241:GLU:N	1.97	0.78
1:A:214:VAL:HG21	1:A:298:LEU:HD22	1.66	0.75
1:B:305:THR:HG22	1:B:307:MET:H	1.56	0.70
1:D:367:ILE:HG13	1:D:368:LEU:HD12	1.72	0.70
1:A:196:SER:HB3	1:A:213:ARG:HB3	1.74	0.70
1:C:247:GLN:NE2	1:D:267:ILE:HD11	2.07	0.69
1:A:203:MET:HE2	1:A:208:LEU:HD13	1.75	0.69
1:A:203:MET:HE2	1:A:208:LEU:CD1	2.22	0.69
1:D:233:LYS:HD2	1:D:292:PHE:HB3	1.73	0.69
1:D:61:VAL:HG23	1:D:217:LEU:HD22	1.74	0.68
1:D:279:LEU:HD22	1:D:282:LEU:HD12	1.76	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:203:MET:HE1	1:B:361:LEU:HD13	1.76	0.67
1:B:223:VAL:HG12	1:B:224:GLU:HG3	1.76	0.67
1:A:239:GLU:HB3	1:B:325:ARG:HH21	1.59	0.67
1:D:58:PHE:HD2	1:D:341:TYR:HE1	1.42	0.67
1:A:216:ASN:HD21	1:A:218:ARG:HB3	1.60	0.67
1:A:214:VAL:CG2	1:A:298:LEU:HD22	2.25	0.66
1:A:325:ARG:HD3	1:D:241:GLU:HG3	1.77	0.66
1:B:301:MET:CG	1:B:308:THR:HG22	2.26	0.65
1:C:223:VAL:HG12	1:C:224:GLU:HG3	1.77	0.65
1:A:188:LYS:HA	1:A:191:GLU:CG	2.26	0.64
1:A:239:GLU:HB3	1:B:325:ARG:NH2	2.14	0.63
1:B:203:MET:HE2	1:B:208:LEU:CD1	2.26	0.62
1:C:196:SER:HB3	1:C:213:ARG:HB3	1.82	0.62
1:B:214:VAL:HG21	1:B:298:LEU:HD22	1.82	0.62
1:D:281:ASP:OD2	1:D:359:ARG:HD2	2.00	0.62
1:A:297:ILE:HD12	1:A:312:ARG:HG2	1.80	0.61
1:D:327:GLU:HG3	1:D:346:LYS:HB3	1.82	0.61
1:A:203:MET:CE	1:A:208:LEU:HD13	2.29	0.61
1:C:247:GLN:HE22	1:D:267:ILE:HD11	1.65	0.61
1:A:267:ILE:HD11	1:D:247:GLN:HE22	1.64	0.60
1:A:216:ASN:HD21	1:A:218:ARG:CB	2.14	0.60
1:B:237:THR:HG23	1:B:239:GLU:N	2.13	0.60
1:D:203:MET:HE2	1:D:208:LEU:HD13	1.84	0.60
1:B:237:THR:CG2	1:B:241:GLU:HB3	2.32	0.60
1:C:214:VAL:HG21	1:C:298:LEU:HD22	1.84	0.59
1:D:297:ILE:CD1	1:D:312:ARG:HG2	2.31	0.59
1:B:261:ILE:H	1:B:261:ILE:HD13	1.67	0.59
1:B:214:VAL:CG2	1:B:298:LEU:HD22	2.33	0.58
1:D:203:MET:CE	1:D:208:LEU:HD13	2.34	0.58
1:D:237:THR:HG22	1:D:238:SER:N	2.19	0.58
1:B:251:ASN:O	1:B:271:HIS:HD2	1.87	0.57
1:B:63:GLU:CD	1:B:63:GLU:H	2.11	0.57
1:B:237:THR:CG2	1:B:241:GLU:H	2.15	0.57
1:B:203:MET:HE3	1:B:354:PRO:HG2	1.86	0.57
1:B:192:THR:HG22	1:B:218:ARG:CZ	2.35	0.56
1:A:307:MET:HE3	1:B:304:ALA:HA	1.86	0.56
1:D:363:GLU:O	1:D:367:ILE:HG23	2.06	0.56
1:A:281:ASP:OD1	1:A:359:ARG:HD2	2.06	0.56
1:C:233:LYS:HD3	1:C:292:PHE:HB3	1.87	0.56
1:D:232:LEU:HG	1:D:295:VAL:HG23	1.88	0.55
1:C:307:MET:HE3	1:D:304:ALA:HA	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:223:VAL:HG12	1:D:224:GLU:HG3	1.89	0.55
1:C:201:ILE:HD13	1:C:210:LEU:HA	1.89	0.55
1:C:214:VAL:HB	1:C:269:ILE:HD11	1.89	0.55
1:B:216:ASN:HB2	1:B:263:LEU:HD13	1.90	0.54
1:D:284:LYS:HD3	1:D:355:LEU:HD22	1.91	0.53
1:A:301:MET:CG	1:A:308:THR:HG22	2.31	0.53
1:C:53:HIS:HB2	2:C:477:HOH:O	2.08	0.53
1:A:253:GLY:HA3	1:A:258:ILE:HG13	1.90	0.53
1:B:232:LEU:HD22	1:B:247:GLN:HG2	1.90	0.53
1:D:203:MET:HE2	1:D:208:LEU:CD1	2.39	0.53
1:D:253:GLY:HA3	1:D:258:ILE:HG13	1.91	0.53
1:D:214:VAL:HG21	1:D:298:LEU:HD22	1.91	0.53
1:A:188:LYS:HG2	1:A:191:GLU:HG3	1.90	0.52
1:A:56:VAL:HA	1:A:337:TYR:O	2.09	0.52
1:A:63:GLU:HG3	1:A:189:ARG:HD2	1.92	0.52
1:D:220:SER:HB2	1:D:303:GLU:HB3	1.92	0.52
1:B:283:SER:H	1:B:286:ASP:HB2	1.74	0.52
1:B:279:LEU:HD22	1:B:282:LEU:HD12	1.91	0.51
1:D:324:HIS:NE2	1:D:349:GLU:HG2	2.25	0.51
1:B:325:ARG:O	1:B:347:THR:HB	2.10	0.51
1:C:192:THR:HB	1:C:218:ARG:HH21	1.76	0.51
1:C:237:THR:HG22	1:C:238:SER:N	2.25	0.51
1:C:220:SER:HB2	1:C:303:GLU:HB3	1.93	0.51
1:B:237:THR:HG21	1:B:241:GLU:HB3	1.92	0.51
1:D:234:SER:HB3	1:D:244:PRO:O	2.11	0.51
1:D:214:VAL:CG1	1:D:269:ILE:HD11	2.41	0.50
1:D:216:ASN:HD21	1:D:218:ARG:HB2	1.77	0.50
1:B:271:HIS:HE1	1:B:277:SER:OG	1.95	0.50
1:C:218:ARG:NH1	1:C:302:VAL:HG21	2.27	0.49
1:B:226:HIS:HD2	2:B:444:HOH:O	1.95	0.49
1:B:301:MET:SD	1:B:308:THR:HG22	2.53	0.49
1:D:201:ILE:HA	1:D:209:CYS:O	2.13	0.48
1:A:203:MET:HE1	1:A:208:LEU:HB2	1.94	0.48
1:B:292:PHE:C	1:B:292:PHE:CD1	2.91	0.48
1:C:232:LEU:CD2	1:C:247:GLN:HG2	2.43	0.48
1:C:303:GLU:O	1:C:304:ALA:HB3	2.13	0.48
1:C:334:LYS:HE3	1:C:335:HIS:NE2	2.28	0.48
1:A:45:SER:N	1:A:335:HIS:HA	2.29	0.48
1:A:283:SER:O	1:A:287:ILE:HG12	2.13	0.48
1:C:292:PHE:CD1	1:C:292:PHE:C	2.93	0.47
1:B:307:MET:HE3	1:C:304:ALA:HA	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:226:HIS:HE1	1:B:224:GLU:OE2	1.97	0.47
1:A:205:ASP:OD2	1:D:237:THR:HG23	2.14	0.47
1:C:301:MET:HG2	1:C:308:THR:HG22	1.97	0.47
1:C:325:ARG:O	1:C:347:THR:HB	2.15	0.47
1:D:214:VAL:CG2	1:D:298:LEU:HD22	2.45	0.47
1:D:216:ASN:HD21	1:D:218:ARG:CB	2.28	0.47
1:D:216:ASN:ND2	1:D:218:ARG:H	2.13	0.47
1:B:245:LEU:HD12	1:C:330:LEU:HD21	1.97	0.46
1:D:232:LEU:HD22	1:D:247:GLN:HG2	1.96	0.46
1:A:49:LYS:HE3	1:A:53:HIS:HB2	1.97	0.46
1:A:237:THR:HG22	1:A:241:GLU:H	1.81	0.46
1:A:203:MET:HE2	1:A:208:LEU:HD12	1.95	0.46
1:D:61:VAL:O	1:D:61:VAL:HG22	2.15	0.46
1:D:58:PHE:CD2	1:D:341:TYR:HE1	2.28	0.46
1:B:216:ASN:HD21	1:B:218:ARG:HB3	1.80	0.46
1:D:272:GLU:HB2	2:D:467:HOH:O	2.14	0.45
1:C:309:THR:HG22	1:C:310:GLN:N	2.30	0.45
1:A:205:ASP:H	1:D:237:THR:CG2	2.30	0.45
1:C:61:VAL:HG23	1:C:217:LEU:HD22	1.99	0.45
1:C:251:ASN:O	1:C:271:HIS:HD2	2.00	0.45
1:D:232:LEU:HD11	1:D:295:VAL:HG21	1.99	0.45
1:A:237:THR:CG2	1:A:241:GLU:H	2.30	0.45
1:A:256:SER:OG	1:A:258:ILE:HG12	2.17	0.45
1:C:214:VAL:CG2	1:C:298:LEU:HD22	2.46	0.45
1:D:277:SER:C	1:D:279:LEU:H	2.25	0.45
1:C:200:VAL:HG12	1:C:321:LEU:HB2	1.99	0.44
1:C:199:ALA:HB3	1:C:320:ILE:CD1	2.48	0.44
1:C:232:LEU:HD22	1:C:247:GLN:HG2	1.99	0.44
1:D:292:PHE:C	1:D:292:PHE:CD1	2.95	0.44
1:C:237:THR:CG2	1:D:205:ASP:H	2.30	0.44
1:A:190:ASN:HB3	1:A:217:LEU:HD13	1.98	0.44
1:A:232:LEU:HD22	1:A:247:GLN:HG2	2.00	0.44
1:A:251:ASN:O	1:A:271:HIS:HD2	2.01	0.44
1:B:284:LYS:HD3	1:B:355:LEU:HD22	2.00	0.44
1:C:226:HIS:HE1	1:D:224:GLU:OE2	2.01	0.44
1:A:190:ASN:HD22	1:A:190:ASN:N	2.15	0.44
1:B:63:GLU:CD	1:B:63:GLU:N	2.76	0.44
1:A:190:ASN:N	1:A:190:ASN:ND2	2.63	0.44
1:A:267:ILE:HD11	1:D:247:GLN:NE2	2.32	0.43
1:B:281:ASP:OD1	1:B:359:ARG:HD2	2.18	0.43
1:D:49:LYS:CE	1:D:53:HIS:HB2	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:58:PHE:CD2	1:A:341:TYR:HE1	2.37	0.43
1:A:241:GLU:HG3	1:B:325:ARG:HD3	2.00	0.43
1:C:250:ILE:HG23	1:C:271:HIS:CD2	2.53	0.43
1:D:228:ARG:HD2	2:D:470:HOH:O	2.19	0.43
1:C:269:ILE:HD12	1:C:269:ILE:H	1.84	0.43
1:A:190:ASN:HB3	1:A:217:LEU:CD1	2.48	0.43
1:D:283:SER:O	1:D:287:ILE:HG12	2.18	0.42
1:B:192:THR:HB	1:B:218:ARG:NH2	2.34	0.42
1:C:212:TRP:HA	1:C:326:TYR:OH	2.20	0.42
1:C:269:ILE:HD12	1:C:269:ILE:N	2.33	0.42
1:C:218:ARG:HE	1:C:218:ARG:HB2	1.74	0.42
1:D:237:THR:CG2	1:D:239:GLU:OE1	2.68	0.42
1:A:237:THR:CG2	1:A:241:GLU:HB3	2.50	0.42
1:B:192:THR:HG22	1:B:218:ARG:NH1	2.35	0.42
1:D:265:SER:O	1:D:266:PRO:C	2.62	0.42
1:C:279:LEU:CD2	1:C:282:LEU:HD12	2.50	0.42
1:C:343:ARG:HG2	1:C:343:ARG:HH11	1.85	0.42
1:A:190:ASN:O	1:A:192:THR:HG23	2.19	0.41
1:A:250:ILE:CD1	1:A:294:ILE:HG23	2.49	0.41
1:A:301:MET:HE2	1:A:301:MET:HB3	1.89	0.41
1:A:237:THR:HG22	1:A:241:GLU:N	2.35	0.41
1:B:309:THR:HG22	1:B:310:GLN:N	2.35	0.41
1:C:253:GLY:HA3	1:C:258:ILE:HG13	2.01	0.41
1:D:49:LYS:HE2	1:D:53:HIS:HB2	2.03	0.41
1:A:334:LYS:C	1:A:336:TYR:N	2.78	0.41
1:C:237:THR:CG2	1:C:239:GLU:OE1	2.68	0.41
1:C:329:VAL:O	1:C:339:VAL:HA	2.21	0.41
1:C:262:PHE:O	1:C:267:ILE:HD11	2.21	0.41
1:A:231:LEU:C	1:A:231:LEU:HD23	2.46	0.41
1:B:324:HIS:CE1	1:B:349:GLU:HG2	2.56	0.41
1:C:201:ILE:HA	1:C:209:CYS:O	2.21	0.41
1:C:226:HIS:HD2	2:C:437:HOH:O	2.03	0.41
1:C:233:LYS:HD3	1:C:292:PHE:CB	2.51	0.41
1:C:192:THR:HB	1:C:218:ARG:NH2	2.36	0.41
1:C:241:GLU:HG3	1:D:325:ARG:HD3	2.03	0.41
1:D:56:VAL:CG2	1:D:339:VAL:HG23	2.51	0.41
1:B:234:SER:OG	1:B:244:PRO:O	2.33	0.40
1:C:216:ASN:ND2	1:C:218:ARG:H	2.19	0.40
1:C:234:SER:OG	1:C:244:PRO:O	2.30	0.40
1:D:237:THR:CG2	1:D:238:SER:N	2.84	0.40
1:C:216:ASN:HD21	1:C:218:ARG:CB	2.35	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	195/270 (72%)	182 (93%)	9 (5%)	4 (2%)	5	6
1	B	191/270 (71%)	176 (92%)	13 (7%)	2 (1%)	12	18
1	C	191/270 (71%)	179 (94%)	12 (6%)	0	100	100
1	D	193/270 (72%)	182 (94%)	11 (6%)	0	100	100
All	All	770/1080 (71%)	719 (93%)	45 (6%)	6 (1%)	16	23

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	335	HIS
1	A	191	GLU
1	A	218	ARG
1	B	244	PRO
1	B	351	PRO
1	A	244	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	177/244 (72%)	173 (98%)	4 (2%)	44	64
1	B	175/244 (72%)	173 (99%)	2 (1%)	65	81

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	175/244 (72%)	173 (99%)	2 (1%)	65	81
1	D	177/244 (72%)	171 (97%)	6 (3%)	32	52
All	All	704/976 (72%)	690 (98%)	14 (2%)	48	68

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	63	GLU
1	A	190	ASN
1	A	263	LEU
1	A	318	ASN
1	B	261	ILE
1	B	263	LEU
1	C	214	VAL
1	C	263	LEU
1	D	55	ASN
1	D	218	ARG
1	D	252	VAL
1	D	263	LEU
1	D	266	PRO
1	D	314	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	190	ASN
1	A	198	ASN
1	A	216	ASN
1	A	226	HIS
1	A	247	GLN
1	A	271	HIS
1	A	318	ASN
1	A	352	ASN
1	B	190	ASN
1	B	197	HIS
1	B	198	ASN
1	B	216	ASN
1	B	221	HIS
1	B	226	HIS
1	B	271	HIS

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Mol	Chain	Res	Type
1	C	60	ASN
1	C	198	ASN
1	C	216	ASN
1	C	221	HIS
1	C	226	HIS
1	C	247	GLN
1	C	271	HIS
1	C	324	HIS
1	C	352	ASN
1	D	57	GLN
1	D	60	ASN
1	D	198	ASN
1	D	216	ASN
1	D	226	HIS
1	D	247	GLN
1	D	271	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	197/270 (72%)	0.49	9 (4%) 37 33	28, 48, 84, 105	0
1	B	195/270 (72%)	0.59	20 (10%) 12 9	26, 46, 77, 100	0
1	C	195/270 (72%)	0.45	12 (6%) 26 23	26, 44, 73, 96	0
1	D	197/270 (72%)	0.64	17 (8%) 16 13	29, 47, 81, 110	0
All	All	784/1080 (72%)	0.54	58 (7%) 20 17	26, 46, 80, 110	0

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	367	ILE	5.3
1	D	368	LEU	5.2
1	D	336	TYR	5.0
1	C	336	TYR	4.2
1	B	304	ALA	3.9
1	D	352	ASN	3.7
1	B	336	TYR	3.6
1	A	45	SER	3.6
1	A	336	TYR	3.4
1	C	360	ASP	3.3
1	B	351	PRO	3.1
1	A	301	MET	2.9
1	B	305	THR	2.9
1	B	349	GLU	2.9
1	B	335	HIS	2.8
1	D	366	TYR	2.8
1	B	192	THR	2.8
1	A	307	MET	2.8
1	B	62	GLY	2.8
1	A	365	LYS	2.6
1	B	365	LYS	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	61	VAL	2.5
1	B	190	ASN	2.5
1	A	291	ASP	2.5
1	B	191	GLU	2.5
1	C	291	ASP	2.4
1	D	361	LEU	2.4
1	D	45	SER	2.4
1	B	301	MET	2.3
1	B	306	ALA	2.3
1	B	307	MET	2.3
1	A	220	SER	2.3
1	D	220	SER	2.3
1	B	362	ALA	2.2
1	A	237	THR	2.2
1	C	364	LYS	2.2
1	D	350	VAL	2.2
1	D	289	ASN	2.2
1	B	356	CYS	2.2
1	C	308	THR	2.2
1	C	365	LYS	2.2
1	C	363	GLU	2.2
1	D	287	ILE	2.1
1	D	188	LYS	2.1
1	B	350	VAL	2.1
1	C	61	VAL	2.1
1	C	245	LEU	2.1
1	B	206	GLY	2.1
1	B	308	THR	2.1
1	D	305	THR	2.1
1	D	61	VAL	2.1
1	C	317	ALA	2.1
1	A	190	ASN	2.1
1	C	287	ILE	2.1
1	C	335	HIS	2.1
1	D	197	HIS	2.1
1	D	192	THR	2.0
1	D	333	GLU	2.0

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

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.