



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 19, 2026 – 07:24 PM UTC

PDB ID : 5KWQ / pdb_00005kwq
Title : Two Tandem RRM Domains of FBP-Interacting Repressor (FIR), also Known as PUF60
Authors : Crichlow, G.V.; Yang, Y.; Zhou, H.; Lolis, E.J.; Braddock, D.T.
Deposited on : 2016-07-18
Resolution : 2.80 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

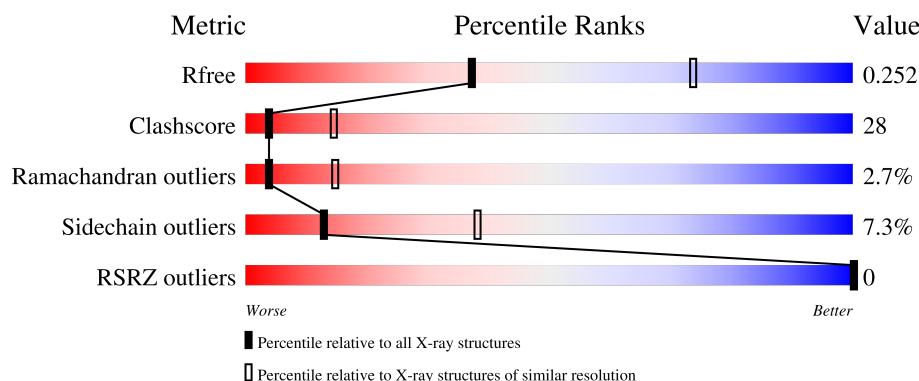
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	216	
1	B	216	

2 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 2896 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 Poly(U)-binding-splicing factor PUF60.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	190	Total	C	N	O	S	0	0	0
			1444	917	243	278	6			
1	B	191	Total	C	N	O	S	0	0	0
			1452	923	244	279	6			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	84	GLY	-	expression tag	UNP Q9UHX1
A	85	SER	-	expression tag	UNP Q9UHX1
A	86	HIS	-	expression tag	UNP Q9UHX1
A	87	MET	-	expression tag	UNP Q9UHX1
A	88	ALA	-	expression tag	UNP Q9UHX1
A	89	SER	-	expression tag	UNP Q9UHX1
A	90	MET	-	expression tag	UNP Q9UHX1
A	91	THR	-	expression tag	UNP Q9UHX1
A	92	GLY	-	expression tag	UNP Q9UHX1
A	93	GLY	-	expression tag	UNP Q9UHX1
A	94	GLN	-	expression tag	UNP Q9UHX1
A	95	GLN	-	expression tag	UNP Q9UHX1
A	96	MET	-	expression tag	UNP Q9UHX1
A	97	GLY	-	expression tag	UNP Q9UHX1
A	98	ARG	-	expression tag	UNP Q9UHX1
A	99	GLY	-	expression tag	UNP Q9UHX1
A	100	SER	-	expression tag	UNP Q9UHX1
A	106	GLY	ARG	engineered mutation	UNP Q9UHX1
A	112	SER	CYS	engineered mutation	UNP Q9UHX1
A	238	ALA	CYS	engineered mutation	UNP Q9UHX1
B	84	GLY	-	expression tag	UNP Q9UHX1
B	85	SER	-	expression tag	UNP Q9UHX1
B	86	HIS	-	expression tag	UNP Q9UHX1
B	87	MET	-	expression tag	UNP Q9UHX1
B	88	ALA	-	expression tag	UNP Q9UHX1

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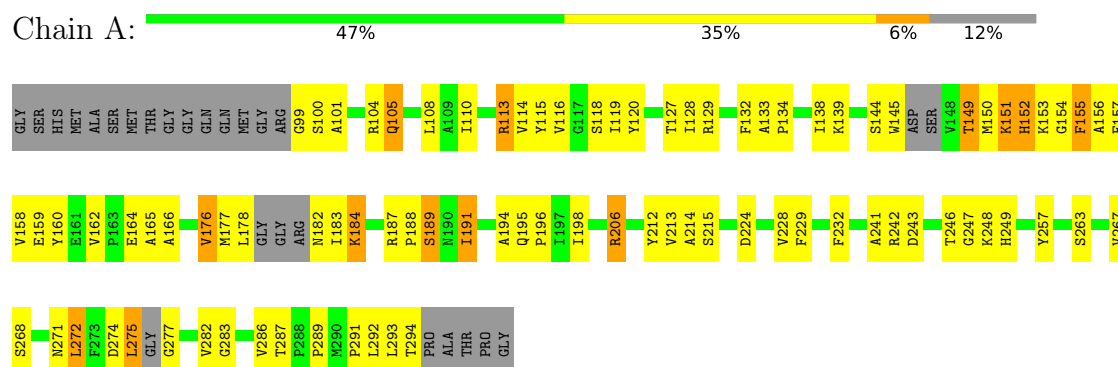
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Chain	Residue	Modelled	Actual	Comment	Reference
B	89	SER	-	expression tag	UNP Q9UHX1
B	90	MET	-	expression tag	UNP Q9UHX1
B	91	THR	-	expression tag	UNP Q9UHX1
B	92	GLY	-	expression tag	UNP Q9UHX1
B	93	GLY	-	expression tag	UNP Q9UHX1
B	94	GLN	-	expression tag	UNP Q9UHX1
B	95	GLN	-	expression tag	UNP Q9UHX1
B	96	MET	-	expression tag	UNP Q9UHX1
B	97	GLY	-	expression tag	UNP Q9UHX1
B	98	ARG	-	expression tag	UNP Q9UHX1
B	99	GLY	-	expression tag	UNP Q9UHX1
B	100	SER	-	expression tag	UNP Q9UHX1
B	106	GLY	ARG	engineered mutation	UNP Q9UHX1
B	112	SER	CYS	engineered mutation	UNP Q9UHX1
B	238	ALA	CYS	engineered mutation	UNP Q9UHX1

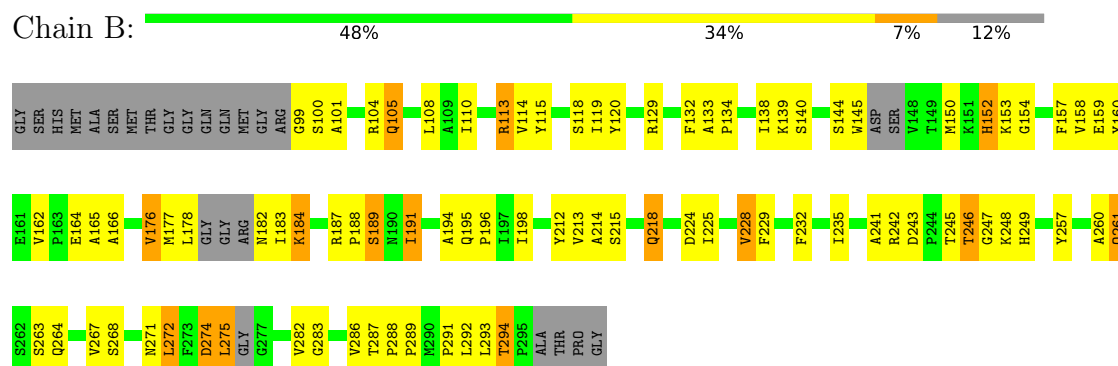
3 Residue-property plots

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: Poly(U)-binding-splicing factor PUF60



- Molecule 1: Poly(U)-binding-splicing factor PUF60



4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α , β , γ	61.83Å 61.83Å 80.41Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.63 – 2.80 19.63 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.7 (19.63-2.80) 99.8 (19.63-2.80)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.42 (at 2.79Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.239 , 0.256 0.236 , 0.252	Depositor DCC
R_{free} test set	408 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å ²)	57.1	Xtriage
Anisotropy	0.045	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 32.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.049 for -h,-k,l 0.487 for h,-h-k,-l 0.055 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2896	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 7.56% 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 [i](#)

5.1 Standard geometry [i](#)

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	0.85	1/1471 (0.1%)	1.16	6/1989 (0.3%)
1	B	0.92	2/1480 (0.1%)	1.18	8/2002 (0.4%)
All	All	0.88	3/2951 (0.1%)	1.17	14/3991 (0.4%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	140	SER	CA-C	-6.46	1.44	1.52
1	A	156	ALA	N-CA	-5.81	1.38	1.45
1	B	140	SER	N-CA	-5.23	1.39	1.45

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	261	GLN	OE1-CD-NE2	-9.79	112.81	122.60
1	B	218	GLN	OE1-CD-NE2	-9.42	113.18	122.60
1	A	155	PHE	CA-CB-CG	6.72	120.52	113.80
1	A	274	ASP	O-C-N	6.56	130.66	122.65
1	B	261	GLN	CG-CD-NE2	6.50	126.15	116.40
1	B	218	GLN	CG-CD-NE2	6.35	125.93	116.40
1	A	151	LYS	CA-C-N	-5.58	110.99	122.82
1	A	151	LYS	C-N-CA	-5.58	110.99	122.82
1	A	152	HIS	N-CA-C	5.56	117.92	109.63
1	A	274	ASP	N-CA-C	5.14	116.74	110.41
1	B	274	ASP	N-CA-C	5.05	116.62	110.41
1	B	228	VAL	CB-CA-C	-5.05	105.33	112.14
1	B	235	ILE	N-CA-C	5.04	115.35	107.99
1	B	152	HIS	O-C-N	5.01	128.76	122.65

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	1444	0	1377	79	0
1	B	1452	0	1386	88	0
All	All	2896	0	2763	161	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 (161) 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:113:ARG:HH11	1:B:113:ARG:HG3	1.01	1.15
1:B:176:VAL:HG13	1:B:177:MET:H	1.20	1.04
1:A:176:VAL:HG13	1:A:177:MET:H	1.20	1.03
1:A:113:ARG:HH11	1:A:113:ARG:HG3	1.28	0.95
1:B:246:THR:HG21	1:B:248:LYS:CG	2.00	0.91
1:A:275:LEU:HA	1:A:277:GLY:N	1.88	0.88
1:B:113:ARG:HG3	1:B:113:ARG:NH1	1.80	0.87
1:B:293:LEU:O	1:B:294:THR:HB	1.76	0.86
1:B:113:ARG:HH11	1:B:113:ARG:CG	1.89	0.83
1:A:177:MET:O	1:A:178:LEU:CB	2.27	0.82
1:A:188:PRO:O	1:A:191:ILE:HG13	1.80	0.81
1:B:246:THR:HG22	1:B:248:LYS:H	1.46	0.81
1:A:120:TYR:HB2	1:A:178:LEU:CB	2.11	0.80
1:B:188:PRO:O	1:B:191:ILE:HG13	1.80	0.80
1:A:263:SER:O	1:A:267:VAL:HG23	1.87	0.75
1:B:177:MET:O	1:B:178:LEU:CB	2.33	0.75
1:A:113:ARG:HG3	1:A:113:ARG:NH1	2.01	0.75
1:A:206:ARG:CB	1:A:206:ARG:HH11	1.99	0.75
1:A:116:VAL:O	1:A:155:PHE:HA	1.87	0.74
1:B:218:GLN:HA	1:B:249:HIS:NE2	2.03	0.74
1:B:263:SER:O	1:B:267:VAL:HG23	1.87	0.74
1:B:246:THR:CG2	1:B:248:LYS:CG	2.65	0.74
1:A:176:VAL:HG13	1:A:177:MET:N	1.99	0.71
1:A:206:ARG:HH11	1:A:206:ARG:HB2	1.56	0.69
1:A:206:ARG:HB2	1:A:206:ARG:NH1	2.07	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:176:VAL:HG13	1:B:177:MET:N	1.99	0.67
1:A:293:LEU:O	1:A:294:THR:HG23	1.95	0.67
1:A:119:ILE:HB	1:A:154:GLY:HA2	1.76	0.66
1:A:246:THR:HG22	1:A:248:LYS:H	1.57	0.66
1:B:274:ASP:O	1:B:275:LEU:HB2	1.96	0.64
1:B:120:TYR:HB2	1:B:178:LEU:CB	2.27	0.64
1:B:194:ALA:O	1:B:198:ILE:HG13	2.00	0.62
1:A:271:ASN:O	1:A:272:LEU:HB2	2.00	0.62
1:A:224:ASP:O	1:A:228:VAL:HG23	2.00	0.62
1:A:194:ALA:O	1:A:198:ILE:HG13	2.00	0.61
1:B:218:GLN:HA	1:B:249:HIS:CE1	2.35	0.61
1:B:271:ASN:O	1:B:272:LEU:HB2	2.00	0.61
1:A:293:LEU:HG	1:A:294:THR:H	1.66	0.60
1:A:99:GLY:O	1:A:104:ARG:HD2	2.01	0.60
1:B:99:GLY:O	1:B:104:ARG:HD2	2.01	0.60
1:A:145:TRP:CD2	1:A:152:HIS:HB3	2.37	0.60
1:B:152:HIS:ND1	1:B:154:GLY:N	2.49	0.60
1:A:286:VAL:HG23	1:A:287:THR:HG23	1.85	0.58
1:B:246:THR:HG22	1:B:247:GLY:N	2.19	0.57
1:A:268:SER:HB2	1:B:268:SER:OG	2.04	0.57
1:A:268:SER:OG	1:B:268:SER:HB2	2.04	0.57
1:B:187:ARG:HB3	1:B:191:ILE:HD11	1.86	0.57
1:B:188:PRO:O	1:B:191:ILE:CG1	2.52	0.57
1:A:187:ARG:HB3	1:A:191:ILE:HD11	1.86	0.57
1:B:218:GLN:HA	1:B:249:HIS:HE2	1.69	0.57
1:A:242:ARG:NH1	1:A:242:ARG:HG2	2.20	0.56
1:A:115:TYR:CE2	1:B:184:LYS:HD3	2.41	0.56
1:B:242:ARG:HG2	1:B:242:ARG:NH1	2.20	0.56
1:A:110:ILE:HA	1:A:113:ARG:NH1	2.20	0.56
1:A:184:LYS:HD3	1:B:115:TYR:CE2	2.41	0.55
1:B:113:ARG:NH1	1:B:113:ARG:CG	2.56	0.55
1:A:188:PRO:O	1:A:191:ILE:CG1	2.52	0.55
1:B:224:ASP:O	1:B:228:VAL:HG23	2.07	0.55
1:A:164:GLU:HG2	1:A:212:TYR:HB2	1.89	0.54
1:B:152:HIS:ND1	1:B:154:GLY:HA2	2.22	0.54
1:B:242:ARG:HG2	1:B:242:ARG:HH11	1.73	0.53
1:B:164:GLU:HG2	1:B:212:TYR:HB2	1.89	0.53
1:B:293:LEU:O	1:B:294:THR:CB	2.51	0.53
1:B:293:LEU:HG	1:B:294:THR:H	1.73	0.53
1:A:242:ARG:HG2	1:A:242:ARG:HH11	1.73	0.52
1:B:162:VAL:HG22	1:B:164:GLU:OE1	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:110:ILE:HA	1:B:113:ARG:NH1	2.24	0.52
1:B:152:HIS:CE1	1:B:154:GLY:HA2	2.45	0.52
1:A:162:VAL:HG22	1:A:164:GLU:OE1	2.10	0.51
1:A:113:ARG:NH1	1:A:113:ARG:CG	2.71	0.51
1:A:241:ALA:O	1:A:249:HIS:HA	2.11	0.51
1:A:104:ARG:O	1:A:108:LEU:HG	2.11	0.51
1:B:104:ARG:O	1:B:108:LEU:HG	2.10	0.51
1:B:241:ALA:O	1:B:249:HIS:HA	2.11	0.50
1:A:206:ARG:CB	1:A:206:ARG:NH1	2.70	0.50
1:A:289:PRO:O	1:A:291:PRO:HD3	2.11	0.50
1:A:176:VAL:O	1:A:177:MET:HB2	2.11	0.50
1:B:132:PHE:HB3	1:B:160:TYR:OH	2.12	0.50
1:A:105:GLN:HB2	1:A:292:LEU:HD22	1.94	0.50
1:B:176:VAL:CG1	1:B:177:MET:H	2.04	0.50
1:A:132:PHE:HB3	1:A:160:TYR:OH	2.12	0.49
1:B:176:VAL:HA	1:B:182:ASN:HB2	1.95	0.49
1:B:145:TRP:HA	1:B:152:HIS:HA	1.94	0.49
1:B:287:THR:HB	1:B:288:PRO:HD2	1.93	0.49
1:B:289:PRO:O	1:B:291:PRO:HD3	2.11	0.49
1:A:242:ARG:HD3	1:A:247:GLY:O	2.13	0.49
1:B:242:ARG:HD3	1:B:247:GLY:O	2.13	0.49
1:B:105:GLN:HB2	1:B:292:LEU:HD22	1.94	0.48
1:A:176:VAL:HA	1:A:182:ASN:HB2	1.95	0.48
1:A:213:VAL:HG22	1:A:282:VAL:HG22	1.96	0.48
1:B:213:VAL:HG22	1:B:282:VAL:HG22	1.96	0.47
1:A:110:ILE:HG22	1:A:191:ILE:HG23	1.96	0.47
1:A:138:ILE:HD12	1:A:158:VAL:CG1	2.45	0.47
1:B:110:ILE:HG22	1:B:191:ILE:HG23	1.96	0.47
1:B:176:VAL:HA	1:B:182:ASN:CB	2.45	0.47
1:B:138:ILE:HD12	1:B:158:VAL:CG1	2.45	0.47
1:B:165:ALA:O	1:B:166:ALA:C	2.58	0.47
1:B:191:ILE:HG22	1:B:198:ILE:CD1	2.45	0.46
1:A:165:ALA:O	1:A:166:ALA:C	2.58	0.46
1:B:118:SER:HB2	1:B:183:ILE:HA	1.98	0.46
1:A:118:SER:HB2	1:A:183:ILE:HA	1.97	0.46
1:A:176:VAL:HA	1:A:182:ASN:CB	2.45	0.46
1:A:176:VAL:HG22	1:A:177:MET:N	2.31	0.46
1:A:243:ASP:O	1:A:247:GLY:N	2.49	0.46
1:B:152:HIS:ND1	1:B:154:GLY:CA	2.78	0.46
1:B:195:GLN:HB3	1:B:196:PRO:HD3	1.97	0.46
1:B:176:VAL:HG22	1:B:177:MET:N	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:282:VAL:HG12	1:A:283:GLY:N	2.31	0.45
1:B:229:PHE:HB3	1:B:257:TYR:OH	2.15	0.45
1:A:144:SER:OG	1:A:153:LYS:CE	2.64	0.45
1:A:191:ILE:HG22	1:A:198:ILE:CD1	2.45	0.45
1:A:195:GLN:HB3	1:A:196:PRO:HD3	1.97	0.45
1:B:229:PHE:O	1:B:232:PHE:HD2	2.00	0.45
1:A:229:PHE:HB3	1:A:257:TYR:OH	2.15	0.45
1:B:282:VAL:HG12	1:B:283:GLY:N	2.31	0.45
1:B:164:GLU:HB3	1:B:212:TYR:CD2	2.52	0.45
1:B:187:ARG:HB3	1:B:191:ILE:CD1	2.47	0.45
1:B:245:THR:O	1:B:245:THR:HG22	2.16	0.45
1:A:164:GLU:HB3	1:A:212:TYR:CD2	2.52	0.44
1:A:212:TYR:CZ	1:A:214:ALA:HB2	2.52	0.44
1:A:115:TYR:CD2	1:B:184:LYS:HD3	2.52	0.44
1:A:271:ASN:O	1:A:272:LEU:CB	2.66	0.44
1:B:212:TYR:CZ	1:B:214:ALA:HB2	2.52	0.44
1:A:242:ARG:NH1	1:A:249:HIS:N	2.66	0.44
1:A:145:TRP:CH2	1:A:152:HIS:HD2	2.35	0.44
1:A:187:ARG:HB3	1:A:191:ILE:CD1	2.47	0.44
1:B:191:ILE:HG13	1:B:191:ILE:H	1.35	0.44
1:B:286:VAL:HG23	1:B:287:THR:HG23	2.00	0.44
1:A:144:SER:OG	1:A:153:LYS:HE2	2.18	0.44
1:A:149:THR:O	1:A:151:LYS:N	2.51	0.44
1:A:229:PHE:O	1:A:232:PHE:HD2	2.00	0.44
1:B:242:ARG:NH1	1:B:249:HIS:N	2.66	0.44
1:A:184:LYS:HD3	1:B:115:TYR:CD2	2.52	0.43
1:A:145:TRP:CE3	1:A:152:HIS:HB3	2.52	0.43
1:B:271:ASN:O	1:B:272:LEU:CB	2.66	0.43
1:A:105:GLN:O	1:A:105:GLN:HG2	2.18	0.43
1:A:133:ALA:N	1:A:134:PRO:CD	2.81	0.43
1:B:184:LYS:H	1:B:184:LYS:HG2	1.70	0.43
1:B:133:ALA:N	1:B:134:PRO:CD	2.81	0.43
1:B:243:ASP:O	1:B:247:GLY:N	2.52	0.43
1:B:293:LEU:HD12	1:B:293:LEU:HA	1.81	0.43
1:B:105:GLN:HG2	1:B:105:GLN:O	2.18	0.42
1:A:184:LYS:H	1:A:184:LYS:HG2	1.70	0.42
1:A:119:ILE:HB	1:A:154:GLY:CA	2.49	0.42
1:B:246:THR:HG22	1:B:248:LYS:CG	2.48	0.42
1:A:157:PHE:CE2	1:A:188:PRO:HG3	2.55	0.42
1:A:100:SER:O	1:A:101:ALA:C	2.62	0.42
1:B:187:ARG:CB	1:B:191:ILE:HD11	2.49	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:260:ALA:O	1:B:261:GLN:C	2.63	0.41
1:B:100:SER:O	1:B:101:ALA:C	2.62	0.41
1:B:157:PHE:CE2	1:B:188:PRO:HG3	2.55	0.41
1:B:119:ILE:O	1:B:120:TYR:C	2.64	0.41
1:A:127:THR:O	1:A:128:ILE:C	2.64	0.41
1:B:224:ASP:O	1:B:225:ILE:C	2.64	0.41
1:B:144:SER:OG	1:B:153:LYS:HD2	2.20	0.41
1:B:245:THR:O	1:B:245:THR:CG2	2.69	0.41
1:A:119:ILE:O	1:A:120:TYR:C	2.63	0.40
1:B:139:LYS:HE2	1:B:159:GLU:OE2	2.22	0.40
1:A:139:LYS:HE2	1:A:159:GLU:OE2	2.22	0.40
1:A:187:ARG:CB	1:A:191:ILE:HD11	2.49	0.40
1:B:176:VAL:CG1	1:B:177:MET:N	2.71	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	182/216 (84%)	167 (92%)	10 (6%)	5 (3%)	4	15
1	B	183/216 (85%)	168 (92%)	10 (6%)	5 (3%)	4	15
All	All	365/432 (84%)	335 (92%)	20 (6%)	10 (3%)	4	15

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	149	THR
1	A	150	MET
1	B	150	MET
1	B	294	THR
1	A	189	SER

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Mol	Chain	Res	Type
1	A	272	LEU
1	B	189	SER
1	B	272	LEU
1	A	176	VAL
1	B	176	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	143/171 (84%)	133 (93%)	10 (7%)	14	40
1	B	144/171 (84%)	133 (92%)	11 (8%)	12	36
All	All	287/342 (84%)	266 (93%)	21 (7%)	13	38

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	105	GLN
1	A	113	ARG
1	A	114	VAL
1	A	129	ARG
1	A	184	LYS
1	A	189	SER
1	A	191	ILE
1	A	206	ARG
1	A	215	SER
1	A	275	LEU
1	B	105	GLN
1	B	113	ARG
1	B	114	VAL
1	B	129	ARG
1	B	184	LYS
1	B	189	SER
1	B	191	ILE
1	B	215	SER

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Mol	Chain	Res	Type
1	B	246	THR
1	B	264	GLN
1	B	275	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	174	ASN
1	B	264	GLN

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 ⓘ

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	190/216 (87%)	-1.54	0 100 100	37, 47, 65, 89	0
1	B	191/216 (88%)	-1.55	0 100 100	34, 47, 66, 97	0
All	All	381/432 (88%)	-1.54	0 100 100	34, 47, 66, 97	0

There are no RSRZ outliers to report.

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.