



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:51 AM GMT

PDB ID : 3JYM
Title : Crystal Structure of the 3 FKBP domains of wheat FKBP73
Authors : Dym, O.; Breiman, A.; Israel Structural Proteomics Center (ISPC)
Deposited on : 2009-09-22
Resolution : 2.28 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

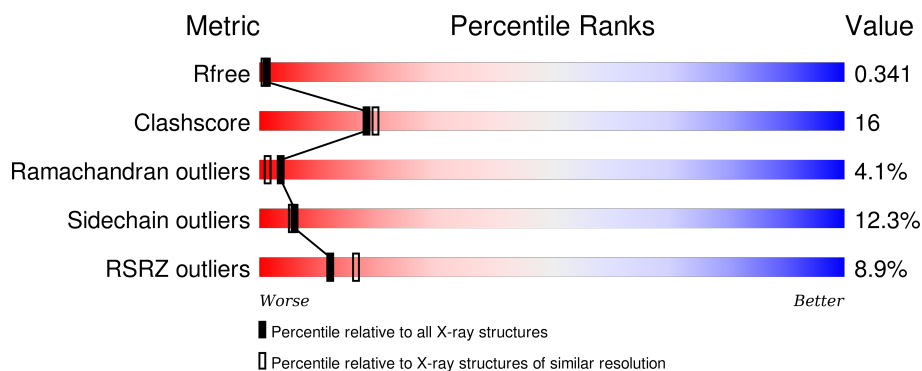
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.28 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}	91344	5193 (2.30-2.26)
Clashscore	102246	5929 (2.30-2.26)
Ramachandran outliers	100387	5851 (2.30-2.26)
Sidechain outliers	100360	5850 (2.30-2.26)
RSRZ outliers	91569	5204 (2.30-2.26)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	377	
1	B	377	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 4726 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 FK506-binding protein (FKBP) from wheat.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	301	Total	C	N	O	S	0	0	0
			2344	1501	378	460	5			
1	B	308	Total	C	N	O	S	0	0	0
			2382	1523	385	469	5			



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	186.96 Å 31.01 Å 68.98 Å 90.00° 93.59° 90.00°	Depositor
Resolution (Å)	46.63 – 2.28 46.65 – 2.28	Depositor EDS
% Data completeness (in resolution range)	98.6 (46.63-2.28) 98.6 (46.65-2.28)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.97 (at 2.27 Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.270 , 0.343 0.271 , 0.341	Depositor DCC
R_{free} test set	1833 reflections (5.25%)	DCC
Wilson B-factor (Å ²)	30.6	Xtriage
Anisotropy	0.167	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 38.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 36734 reflections	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	4726	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.84% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	1.68	32/2382 (1.3%)	1.29	10/3210 (0.3%)
1	B	1.54	19/2419 (0.8%)	1.25	11/3261 (0.3%)
All	All	1.61	51/4801 (1.1%)	1.27	21/6471 (0.3%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	1	1
1	B	0	4
All	All	1	5

All (51) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	209	CYS	CB-SG	-12.13	1.61	1.82
1	A	209	CYS	CB-SG	-11.82	1.62	1.82
1	A	355	TYR	CD2-CE2	-7.48	1.28	1.39
1	A	244	GLY	N-CA	7.44	1.57	1.46
1	A	283	GLU	CB-CG	-7.09	1.38	1.52
1	A	355	TYR	CD1-CE1	-6.50	1.29	1.39
1	B	379	GLU	CD-OE1	-6.50	1.18	1.25
1	A	273	ASP	CB-CG	-6.15	1.38	1.51
1	A	377	GLU	CD-OE1	-6.14	1.18	1.25
1	B	201	PHE	CE1-CZ	6.00	1.48	1.37
1	B	166	GLU	CG-CD	5.99	1.60	1.51
1	A	268	THR	CB-CG2	-5.95	1.32	1.52
1	A	199	VAL	CB-CG2	5.93	1.65	1.52
1	A	377	GLU	CD-OE2	-5.91	1.19	1.25
1	B	200	GLU	CD-OE2	-5.86	1.19	1.25
1	B	194	SER	CB-OG	-5.85	1.34	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	213	ALA	CA-CB	-5.70	1.40	1.52
1	A	313	LYS	CE-NZ	-5.59	1.35	1.49
1	A	213	ALA	CA-CB	-5.56	1.40	1.52
1	A	343	LYS	CE-NZ	5.54	1.62	1.49
1	B	371	ASN	CB-CG	5.50	1.63	1.51
1	A	263	SER	CB-OG	-5.47	1.35	1.42
1	A	357	TYR	CB-CG	-5.46	1.43	1.51
1	B	247	GLY	N-CA	5.46	1.54	1.46
1	A	296	VAL	CB-CG1	-5.42	1.41	1.52
1	B	265	LYS	CB-CG	5.41	1.67	1.52
1	B	183	TYR	CD2-CE2	5.39	1.47	1.39
1	A	312	LYS	C-O	-5.36	1.13	1.23
1	A	165	LYS	CE-NZ	5.36	1.62	1.49
1	A	337	ALA	C-O	-5.35	1.13	1.23
1	B	262	VAL	CB-CG2	-5.33	1.41	1.52
1	A	371	ASN	CB-CG	5.27	1.63	1.51
1	A	312	LYS	CB-CG	-5.26	1.38	1.52
1	A	283	GLU	CG-CD	-5.26	1.44	1.51
1	B	301	THR	CB-CG2	-5.23	1.35	1.52
1	A	37	GLU	C-O	5.22	1.33	1.23
1	A	196	SER	CB-OG	-5.22	1.35	1.42
1	A	292	GLU	CG-CD	5.20	1.59	1.51
1	B	379	GLU	CD-OE2	-5.17	1.20	1.25
1	A	176	PRO	CA-C	5.13	1.63	1.52
1	B	283	GLU	CD-OE2	5.13	1.31	1.25
1	A	379	GLU	CD-OE1	-5.09	1.20	1.25
1	B	160	PHE	CD1-CE1	5.09	1.49	1.39
1	A	181	VAL	CA-CB	5.08	1.65	1.54
1	A	194	SER	CB-OG	-5.06	1.35	1.42
1	A	379	GLU	CD-OE2	-5.04	1.20	1.25
1	B	297	THR	CB-CG2	-5.04	1.35	1.52
1	A	319	GLU	C-O	-5.04	1.13	1.23
1	A	164	LEU	N-CA	-5.04	1.36	1.46
1	B	356	ALA	CA-CB	5.03	1.63	1.52
1	B	256	VAL	CB-CG2	5.01	1.63	1.52

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	247	GLY	N-CA-C	-8.12	92.80	113.10
1	A	240	ARG	NE-CZ-NH1	-7.66	116.47	120.30
1	B	316	ASP	CB-CG-OD1	-7.08	111.93	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	244	GLY	N-CA-C	-6.96	95.69	113.10
1	B	88	LYS	N-CA-C	-6.31	93.97	111.00
1	B	306	ASP	CB-CG-OD1	6.20	123.88	118.30
1	A	240	ARG	NE-CZ-NH2	6.18	123.39	120.30
1	B	156	ASP	CB-CG-OD1	-6.16	112.76	118.30
1	B	227	LEU	CB-CG-CD1	-5.84	101.06	111.00
1	A	286	GLY	N-CA-C	5.84	127.71	113.10
1	A	227	LEU	CB-CG-CD1	-5.81	101.12	111.00
1	B	273	ASP	CB-CA-C	-5.69	99.02	110.40
1	B	287	TYR	CB-CA-C	5.61	121.62	110.40
1	A	177	ASP	CB-CG-OD2	5.53	123.28	118.30
1	B	234	GLY	N-CA-C	-5.19	100.12	113.10
1	A	273	ASP	N-CA-C	5.18	124.99	111.00
1	A	268	THR	OG1-CB-CG2	5.17	121.90	110.00
1	A	244	GLY	N-CA-C	-5.16	100.19	113.10
1	A	289	ARG	NE-CZ-NH1	5.14	122.87	120.30
1	B	240	ARG	NE-CZ-NH1	-5.09	117.75	120.30
1	B	145	LEU	CA-CB-CG	5.04	126.88	115.30

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	268	THR	CB

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	245	GLU	Peptide
1	B	155	LYS	Peptide
1	B	243	ALA	Peptide
1	B	286	GLY	Peptide
1	B	87	PHE	Peptide

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	2344	0	2347	63	0
1	B	2382	0	2392	85	0
All	All	4726	0	4739	148	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 (148) 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:273:ASP:OD1	1:A:275:LYS:HD2	1.27	1.33
1:B:273:ASP:OD1	1:B:275:LYS:CD	1.90	1.18
1:B:273:ASP:OD1	1:B:275:LYS:HD2	0.93	1.10
1:B:243:ALA:HB1	1:B:244:GLY:HA3	1.11	1.06
1:A:65:VAL:O	1:A:87:PHE:N	2.00	0.94
1:B:35:GLU:OE1	1:B:113:ALA:HB1	1.67	0.94
1:B:243:ALA:CB	1:B:244:GLY:HA3	1.98	0.94
1:B:243:ALA:HB1	1:B:244:GLY:CA	1.99	0.91
1:B:324:LYS:HD3	1:B:327:GLU:OE2	1.74	0.87
1:A:175:ASP:OD1	1:A:217:LYS:HE3	1.77	0.84
1:B:65:VAL:H	1:B:87:PHE:N	1.77	0.82
1:A:269:GLU:HB3	1:A:274:LYS:HB3	1.64	0.79
1:B:183:TYR:OH	1:B:207:HIS:HE1	1.67	0.78
1:A:273:ASP:OD1	1:A:275:LYS:CD	2.23	0.73
1:B:311:LEU:HD12	1:B:312:LYS:N	2.05	0.72
1:A:89:PHE:HE1	1:A:101:ASP:OD1	1.74	0.69
1:A:87:PHE:HA	1:A:88:LYS:HB2	1.75	0.69
1:B:134:ALA:O	1:B:136:ALA:N	2.25	0.69
1:B:108:LYS:O	1:B:111:GLU:HB3	1.93	0.69
1:B:65:VAL:O	1:B:87:PHE:HB2	1.93	0.68
1:A:184:GLU:OE2	1:A:186:ARG:NH1	2.27	0.68
1:A:63:VAL:CG1	1:A:91:LEU:HD12	2.25	0.66
1:B:107:MET:HE3	1:B:111:GLU:HG2	1.78	0.64
1:A:34:GLU:HG2	1:A:35:GLU:H	1.62	0.64
1:B:226:LEU:C	1:B:226:LEU:HD23	2.18	0.64
1:B:152:ASP:HB2	1:B:160:PHE:CE1	2.33	0.64
1:A:90:LYS:HB3	1:A:93:GLN:HG3	1.80	0.63
1:A:226:LEU:HD21	1:A:256:VAL:HG13	1.80	0.63
1:B:107:MET:HE1	1:B:142:VAL:O	1.99	0.62
1:B:117:ILE:O	1:B:117:ILE:CG2	2.48	0.61
1:A:367:ILE:O	1:A:369:PRO:HD3	2.02	0.60
1:A:237:GLU:O	1:A:251:PRO:HB3	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:352:PRO:HA	1:B:373:THR:HB	1.84	0.59
1:B:293:GLY:HA2	1:B:324:LYS:HE3	1.84	0.58
1:A:63:VAL:HG11	1:A:91:LEU:HD12	1.85	0.58
1:B:237:GLU:HA	1:B:252:ASN:OD1	2.03	0.58
1:B:38:ILE:O	1:B:112:ASN:HB2	2.03	0.58
1:B:34:GLU:CG	1:B:35:GLU:H	2.17	0.57
1:A:183:TYR:HA	1:A:258:ASP:O	2.03	0.57
1:B:154:ALA:C	1:B:156:ASP:HB2	2.25	0.57
1:B:367:ILE:O	1:B:369:PRO:HD3	2.05	0.57
1:A:97:ILE:O	1:A:99:GLY:N	2.37	0.57
1:A:179:VAL:O	1:A:200:GLU:HA	2.06	0.56
1:B:91:LEU:HD21	1:B:105:LYS:HB2	1.86	0.56
1:B:243:ALA:CB	1:B:244:GLY:CA	2.70	0.56
1:B:38:ILE:HB	1:B:112:ASN:HB3	1.86	0.56
1:A:311:LEU:HD12	1:A:312:LYS:N	2.20	0.56
1:B:200:GLU:OE2	1:B:265:LYS:NZ	2.35	0.55
1:A:208:LEU:HB3	1:A:212:LEU:HD22	1.88	0.55
1:B:293:GLY:O	1:B:324:LYS:HE3	2.07	0.55
1:B:111:GLU:HG3	1:B:112:ASN:O	2.07	0.55
1:A:273:ASP:HB3	1:A:275:LYS:H	1.71	0.54
1:B:293:GLY:O	1:B:324:LYS:HG3	2.08	0.54
1:A:34:GLU:CG	1:A:35:GLU:H	2.19	0.54
1:A:111:GLU:HG2	1:A:112:ASN:H	1.72	0.54
1:A:305:GLN:HG2	1:A:375:ILE:HD11	1.90	0.54
1:A:312:LYS:HG2	1:A:315:HIS:CE1	2.43	0.54
1:B:117:ILE:HG22	1:B:117:ILE:O	2.07	0.53
1:B:34:GLU:HG2	1:B:35:GLU:H	1.74	0.53
1:B:183:TYR:HA	1:B:258:ASP:O	2.09	0.52
1:A:111:GLU:HG2	1:A:112:ASN:N	2.24	0.52
1:A:291:ASN:HB3	1:A:385:LYS:HE3	1.91	0.52
1:B:156:ASP:HB3	1:B:158:GLY:H	1.74	0.52
1:A:33:GLY:HA3	1:A:118:PRO:HD3	1.91	0.52
1:A:235:PHE:CE1	1:A:240:ARG:HD3	2.45	0.51
1:A:230:LYS:HB3	1:A:231:PRO:HD2	1.93	0.51
1:A:67:TYR:HA	1:A:141:ASP:O	2.10	0.51
1:B:63:VAL:HG11	1:B:91:LEU:HD12	1.93	0.51
1:B:330:VAL:HB	1:B:334:LEU:HD23	1.91	0.51
1:B:208:LEU:HB3	1:B:212:LEU:HD22	1.91	0.51
1:B:157:GLY:HA2	1:B:160:PHE:CZ	2.45	0.50
1:B:35:GLU:HA	1:B:114:LEU:O	2.12	0.50
1:B:33:GLY:HA3	1:B:116:THR:O	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:133:PRO:O	1:B:134:ALA:O	2.29	0.50
1:B:152:ASP:HB2	1:B:160:PHE:HE1	1.76	0.50
1:B:224:LYS:HG2	1:B:260:GLU:HG3	1.93	0.50
1:B:297:THR:HG23	1:B:382:SER:OG	2.11	0.49
1:B:57:PRO:HG3	1:B:147:TRP:CD2	2.48	0.49
1:A:109:LYS:O	1:A:144:LEU:O	2.31	0.49
1:B:101:ASP:O	1:B:105:LYS:HB3	2.13	0.48
1:B:35:GLU:OE1	1:B:113:ALA:CB	2.51	0.48
1:B:175:ASP:OD1	1:B:217:LYS:HE3	2.14	0.48
1:B:293:GLY:CA	1:B:324:LYS:HE3	2.43	0.48
1:A:62:GLU:HG3	1:A:90:LYS:HG3	1.95	0.48
1:B:354:GLU:OE2	1:B:354:GLU:N	2.40	0.48
1:B:65:VAL:N	1:B:87:PHE:N	2.57	0.47
1:A:312:LYS:HG2	1:A:315:HIS:NE2	2.29	0.47
1:A:278:LYS:HG3	1:A:349:VAL:HG22	1.95	0.47
1:B:330:VAL:CB	1:B:334:LEU:HD23	2.44	0.47
1:B:293:GLY:O	1:B:324:LYS:CG	2.62	0.47
1:A:63:VAL:HG13	1:A:91:LEU:HD12	1.93	0.47
1:B:57:PRO:HG3	1:B:147:TRP:CE3	2.49	0.47
1:A:54:TRP:CE3	1:A:54:TRP:HA	2.50	0.47
1:A:278:LYS:HE3	1:A:278:LYS:HB3	1.75	0.46
1:A:168:ASP:C	1:A:169:LYS:HG3	2.36	0.46
1:A:285:GLU:OE2	1:A:343:LYS:HD3	2.14	0.46
1:A:352:PRO:HA	1:A:373:THR:HB	1.97	0.46
1:A:54:TRP:HE3	1:A:108:LYS:HZ1	1.62	0.46
1:B:362:SER:HB3	1:B:368:VAL:HB	1.98	0.46
1:A:64:GLU:HG3	1:A:88:LYS:HD2	1.98	0.46
1:A:361:GLU:OE2	1:A:363:LYS:NZ	2.44	0.45
1:A:270:ILE:HG21	1:A:270:ILE:HD13	1.75	0.45
1:A:281:LEU:HD11	1:A:348:LEU:HG	1.97	0.45
1:A:270:ILE:N	1:A:276:ILE:O	2.32	0.45
1:A:54:TRP:HA	1:A:54:TRP:HE3	1.81	0.45
1:B:286:GLY:HA2	1:B:287:TYR:HB3	1.99	0.44
1:B:270:ILE:HG21	1:B:270:ILE:HD13	1.60	0.44
1:B:179:VAL:O	1:B:200:GLU:HA	2.18	0.44
1:B:176:PRO:HA	1:B:204:LYS:HE2	1.99	0.44
1:A:279:LYS:HE2	1:A:279:LYS:HB3	1.84	0.44
1:A:87:PHE:CA	1:A:88:LYS:HB2	2.46	0.44
1:A:170:TRP:HA	1:A:220:LYS:HE3	1.98	0.44
1:B:171:GLU:HG3	1:B:264:TRP:CZ2	2.52	0.44
1:B:57:PRO:HB2	1:B:91:LEU:HD13	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:296:VAL:HG13	1:A:298:VAL:HG13	2.00	0.44
1:A:114:LEU:HD11	1:A:139:GLN:HG2	1.99	0.43
1:B:237:GLU:O	1:B:251:PRO:HB3	2.19	0.43
1:B:286:GLY:CA	1:B:287:TYR:HB3	2.48	0.43
1:B:168:ASP:HB3	1:B:169:LYS:HE2	2.00	0.43
1:B:103:GLY:O	1:B:107:MET:HG3	2.18	0.43
1:B:56:THR:HA	1:B:57:PRO:HD3	1.91	0.43
1:A:224:LYS:HG2	1:A:260:GLU:HG3	2.00	0.43
1:A:117:ILE:O	1:A:138:LEU:N	2.52	0.43
1:B:324:LYS:CD	1:B:327:GLU:OE2	2.58	0.43
1:B:330:VAL:HG21	1:B:334:LEU:HD23	2.00	0.43
1:B:297:THR:CG2	1:B:382:SER:OG	2.67	0.43
1:B:163:ILE:HD13	1:B:163:ILE:HG21	1.62	0.42
1:A:67:TYR:CD1	1:A:68:THR:N	2.88	0.42
1:A:154:ALA:O	1:A:155:LYS:HB2	2.20	0.42
1:A:291:ASN:CB	1:A:385:LYS:HE3	2.49	0.42
1:A:212:LEU:HD12	1:A:212:LEU:HA	1.82	0.42
1:A:250:PRO:O	1:A:251:PRO:C	2.58	0.42
1:A:309:VAL:HG11	1:A:312:LYS:HE2	2.01	0.42
1:B:183:TYR:OH	1:B:207:HIS:CE1	2.59	0.41
1:B:310:PHE:O	1:B:364:GLN:CB	2.68	0.41
1:B:317:GLU:O	1:B:317:GLU:HG2	2.18	0.41
1:B:296:VAL:HG13	1:B:298:VAL:HG13	2.02	0.41
1:B:37:GLU:HG3	1:B:38:ILE:N	2.35	0.41
1:A:156:ASP:HB3	1:A:158:GLY:H	1.85	0.41
1:B:267:VAL:HA	1:B:278:LYS:O	2.20	0.41
1:B:261:LEU:HD12	1:B:261:LEU:HA	1.80	0.41
1:B:107:MET:CE	1:B:111:GLU:HG2	2.47	0.41
1:B:330:VAL:CG2	1:B:334:LEU:HD23	2.51	0.41
1:B:289:ARG:NH1	1:B:340:ASN:O	2.53	0.41
1:B:273:ASP:HB3	1:B:275:LYS:H	1.86	0.41
1:B:208:LEU:HA	1:B:208:LEU:HD12	1.86	0.40
1:A:117:ILE:HA	1:A:118:PRO:HD3	1.95	0.40
1:B:134:ALA:C	1:B:136:ALA:N	2.75	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	292/377 (78%)	261 (89%)	17 (6%)	14 (5%)	3	1
1	B	299/377 (79%)	266 (89%)	23 (8%)	10 (3%)	5	2
All	All	591/754 (78%)	527 (89%)	40 (7%)	24 (4%)	3	1

All (24) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	109	LYS
1	A	246	GLY
1	B	134	ALA
1	B	135	ASN
1	B	287	TYR
1	A	59	VAL
1	A	88	LYS
1	A	98	LYS
1	A	99	GLY
1	A	251	PRO
1	A	286	GLY
1	B	146	SER
1	B	238	MET
1	A	110	GLY
1	B	155	LYS
1	B	251	PRO
1	A	97	ILE
1	A	208	LEU
1	B	243	ALA
1	A	169	LYS
1	A	239	GLY
1	A	273	ASP
1	B	239	GLY
1	B	246	GLY

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	254/311 (82%)	226 (89%)	28 (11%)	8	8
1	B	258/311 (83%)	223 (86%)	35 (14%)	5	4
All	All	512/622 (82%)	449 (88%)	63 (12%)	6	5

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	GLU
1	A	54	TRP
1	A	56	THR
1	A	68	THR
1	A	87	PHE
1	A	91	LEU
1	A	109	LYS
1	A	117	ILE
1	A	149	SER
1	A	156	ASP
1	A	169	LYS
1	A	182	LYS
1	A	197	GLU
1	A	204	LYS
1	A	212	LEU
1	A	217	LYS
1	A	232	GLN
1	A	238	MET
1	A	273	ASP
1	A	283	GLU
1	A	287	TYR
1	A	297	THR
1	A	301	THR
1	A	308	THR
1	A	313	LYS
1	A	360	THR
1	A	363	LYS

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Mol	Chain	Res	Type
1	A	373	THR
1	B	38	ILE
1	B	56	THR
1	B	66	HIS
1	B	91	LEU
1	B	93	GLN
1	B	105	LYS
1	B	111	GLU
1	B	114	LEU
1	B	117	ILE
1	B	135	ASN
1	B	138	LEU
1	B	148	THR
1	B	156	ASP
1	B	166	GLU
1	B	171	GLU
1	B	194	SER
1	B	212	LEU
1	B	226	LEU
1	B	232	GLN
1	B	238	MET
1	B	245	GLU
1	B	251	PRO
1	B	268	THR
1	B	273	ASP
1	B	287	TYR
1	B	296	VAL
1	B	297	THR
1	B	301	THR
1	B	311	LEU
1	B	328	GLU
1	B	348	LEU
1	B	360	THR
1	B	363	LYS
1	B	367	ILE
1	B	373	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	HIS
1	A	232	GLN

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Mol	Chain	Res	Type
1	B	207	HIS
1	B	232	GLN

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 carbohydrates 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	300/377 (79%)	0.58	30 (10%) 9 13	8, 26, 67, 81	0
1	B	307/377 (81%)	0.47	24 (7%) 16 21	11, 30, 61, 80	0
All	All	607/754 (80%)	0.53	54 (8%) 12 16	8, 28, 63, 81	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	100	TRP	7.6
1	A	59	VAL	6.7
1	A	54	TRP	6.5
1	B	59	VAL	6.2
1	B	96	VAL	4.9
1	A	56	THR	4.6
1	B	115	PHE	4.4
1	B	97	ILE	4.4
1	B	100	TRP	4.2
1	A	35	GLU	4.1
1	A	104	ILE	3.5
1	A	144	LEU	3.4
1	A	145	LEU	3.4
1	A	94	GLY	3.4
1	A	96	VAL	3.3
1	A	112	ASN	3.3
1	A	95	GLN	3.3
1	A	115	PHE	3.1
1	A	36	ASN	3.0
1	A	93	GLN	3.0
1	A	170	TRP	3.0
1	B	65	VAL	2.9
1	A	146	SER	2.9
1	B	95	GLN	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	133	PRO	2.8
1	B	38	ILE	2.7
1	B	113	ALA	2.7
1	B	57	PRO	2.7
1	A	65	VAL	2.6
1	B	132	ILE	2.5
1	B	112	ASN	2.5
1	B	69	GLY	2.5
1	B	116	THR	2.4
1	A	57	PRO	2.4
1	B	104	ILE	2.4
1	A	99	GLY	2.4
1	B	366	ALA	2.4
1	A	286	GLY	2.2
1	B	139	GLN	2.2
1	A	105	LYS	2.2
1	B	36	ASN	2.2
1	A	117	ILE	2.2
1	A	138	LEU	2.2
1	A	141	ASP	2.1
1	A	92	GLY	2.1
1	A	97	ILE	2.1
1	B	155	LYS	2.1
1	B	141	ASP	2.1
1	B	55	ASP	2.1
1	A	66	HIS	2.1
1	B	145	LEU	2.0
1	B	286	GLY	2.0
1	A	113	ALA	2.0
1	A	114	LEU	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 ⓘ

There are no carbohydrates 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.