



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 26, 2014 – 03:02 PM GMT

PDB ID : 3GC3
Title : Crystal Structure of Arrestin2S and Clathrin
Authors : Williams, J.C.; Kang, D.S.
Deposited on : 2009-02-21
Resolution : 2.20 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

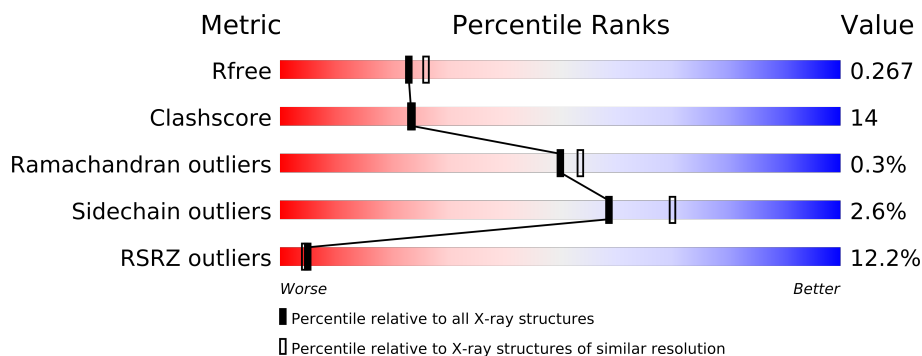
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance



The reported resolution of this entry is 2.20 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}	66092	2938 (2.20-2.20)
Clashscore	79885	3751 (2.20-2.20)
Ramachandran outliers	78287	3681 (2.20-2.20)
Sidechain outliers	78261	3682 (2.20-2.20)
RSRZ outliers	66119	2939 (2.20-2.20)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	385	
2	B	363	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6221 atoms, of which 634 are hydrogens and 0 are deuterium.

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 Beta-arrestin-1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	329	Total	C	H	N	O	S	25	0	0
			2904	1660	322	442	470	10			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	LEU	VARIANT	UNP P17870
A	?	-	LEU	VARIANT	UNP P17870
A	?	-	GLY	VARIANT	UNP P17870
A	?	-	ASP	VARIANT	UNP P17870
A	?	-	LEU	VARIANT	UNP P17870
A	?	-	ALA	VARIANT	UNP P17870
A	?	-	SER	VARIANT	UNP P17870
A	?	-	SER	?	UNP P17870

- Molecule 2 is a protein called Clathrin heavy chain 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	349	Total	C	H	N	O	S	0	0	0
			3041	1738	312	469	504	18			

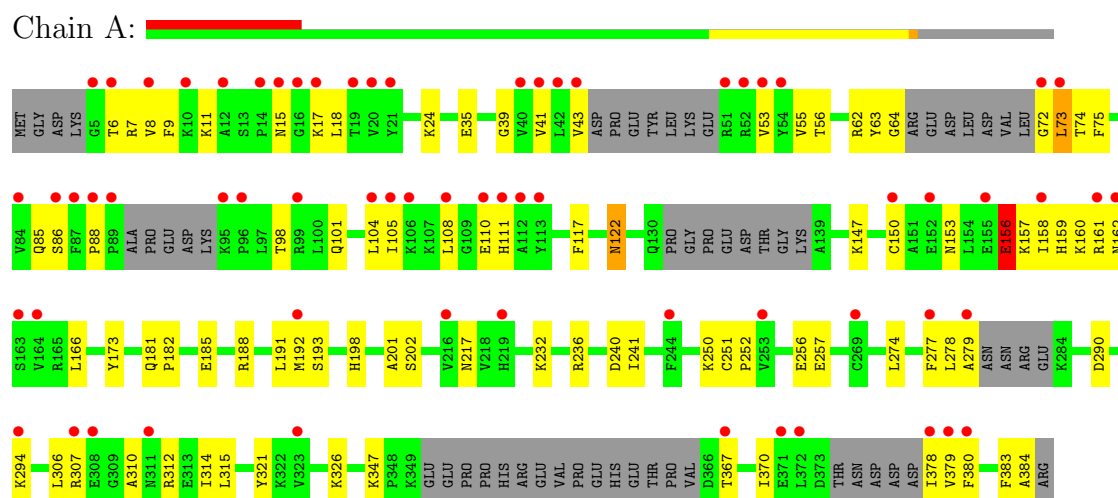
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	133	Total	O	0	0
			133	133		
3	B	143	Total	O	0	0
			143	143		

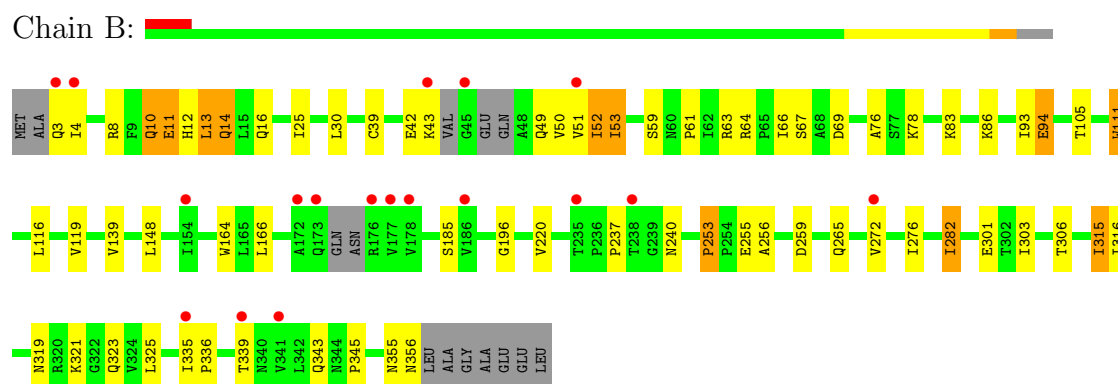
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: Beta-arrestin-1



• Molecule 2: Clathrin heavy chain 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	72.86Å 126.17Å 129.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.07 – 2.20 35.06 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.7 (35.07-2.20) 99.7 (35.06-2.20)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.89 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.4.0034	Depositor
R, R_{free}	0.200 , 0.251 0.229 , 0.267	Depositor DCC
R_{free} test set	3098 reflections (5.07%)	DCC
Wilson B-factor (Å ²)	36.4	Xtriage
Anisotropy	0.044	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 30.2	EDS
Estimated twinning fraction	0.011 for -h,l,k	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 61118 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6221	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.48% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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	1.01	4/2631 (0.2%)	1.00	4/3560 (0.1%)
2	B	1.11	8/2784 (0.3%)	0.99	3/3770 (0.1%)
All	All	1.06	12/5415 (0.2%)	1.00	7/7330 (0.1%)

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	0	2
2	B	0	1
All	All	0	3

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	39	CYS	CB-SG	-8.00	1.68	1.82
2	B	272	VAL	CB-CG2	-7.43	1.37	1.52
2	B	139	VAL	CB-CG2	7.09	1.67	1.52
2	B	119	VAL	CB-CG1	-6.96	1.38	1.52
1	A	156	GLU	CG-CD	-6.48	1.42	1.51
1	A	153	ASN	CB-CG	-5.85	1.37	1.51
1	A	122	ASN	CG-OD1	5.68	1.36	1.24
2	B	11	GLU	CB-CG	5.67	1.62	1.52
2	B	111	TRP	CB-CG	5.61	1.60	1.50
2	B	42	GLU	CB-CG	5.29	1.62	1.52
1	A	321	TYR	CD1-CE1	5.20	1.47	1.39
2	B	10	GLN	CD-OE1	5.12	1.35	1.24

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	188	ARG	NE-CZ-NH2	-13.82	113.39	120.30
1	A	188	ARG	NE-CZ-NH1	11.35	125.98	120.30
2	B	166	LEU	CA-CB-CG	-8.36	96.08	115.30
1	A	188	ARG	CG-CD-NE	-6.27	98.63	111.80
1	A	240	ASP	CB-CG-OD1	5.33	123.10	118.30
2	B	13	LEU	CA-CB-CG	5.15	127.14	115.30
2	B	259	ASP	CB-CG-OD2	-5.14	113.67	118.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	156	GLU	Sidechain
1	A	290	ASP	Peptide
2	B	253	PRO	Peptide

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2582	322	2305	91	0
2	B	2729	312	2428	64	0
3	A	133	0	0	8	0
3	B	143	0	0	4	0
All	All	5587	634	4733	154	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 14.

All (154) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:256:GLU:CG	1:A:274:LEU:HD21	1.39	1.51
1:A:147:LYS:NZ	1:A:158:ILE:HD13	1.27	1.45
1:A:122:ASN:HB2	3:A:509:HOH:O	1.42	1.18
1:A:256:GLU:HG2	1:A:274:LEU:CD2	1.71	1.18
1:A:147:LYS:NZ	1:A:158:ILE:CD1	2.06	1.18

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:8:ARG:NH1	2:B:10:GLN:NE2	1.97	1.12
1:A:147:LYS:HZ1	1:A:158:ILE:CD1	1.61	1.11
1:A:256:GLU:CG	1:A:274:LEU:CD2	2.27	1.09
1:A:294:LYS:HE3	3:A:415:HOH:O	1.53	1.07
1:A:256:GLU:HG3	1:A:274:LEU:HD21	1.39	1.01
2:B:265:GLN:NE2	3:B:410:HOH:O	1.95	0.98
2:B:253:PRO:HD2	2:B:256:ALA:HB3	1.49	0.93
2:B:8:ARG:HH12	2:B:10:GLN:CD	1.73	0.90
2:B:8:ARG:NH1	2:B:10:GLN:CD	2.24	0.90
2:B:8:ARG:HH11	2:B:10:GLN:NE2	1.64	0.89
2:B:335:ILE:H	2:B:335:ILE:HD12	1.34	0.89
1:A:147:LYS:HZ2	1:A:158:ILE:HD13	1.38	0.89
1:A:256:GLU:HG2	1:A:274:LEU:HD21	0.85	0.84
1:A:15:ASN:ND2	1:A:17:LYS:CG	2.44	0.81
2:B:315:ILE:HD12	2:B:316:ILE:N	1.96	0.80
2:B:306:THR:HB	2:B:315:ILE:HD11	1.63	0.80
1:A:15:ASN:ND2	1:A:17:LYS:HG2	1.97	0.79
2:B:355:ASN:O	2:B:356:ASN:HB2	1.82	0.79
1:A:147:LYS:HZ1	1:A:158:ILE:HD13	0.99	0.79
2:B:8:ARG:HH11	2:B:10:GLN:HE21	1.31	0.78
1:A:147:LYS:HZ3	1:A:158:ILE:CD1	1.97	0.78
1:A:256:GLU:HG3	1:A:274:LEU:CD2	2.03	0.77
1:A:147:LYS:HZ3	1:A:158:ILE:HD13	1.50	0.76
1:A:159:HIS:O	1:A:161:ARG:N	2.19	0.76
1:A:63:TYR:O	1:A:73:LEU:CD2	2.36	0.74
2:B:253:PRO:HB3	2:B:255:GLU:OE1	1.88	0.73
1:A:278:LEU:O	1:A:279:ALA:C	2.26	0.73
2:B:76:ALA:CB	2:B:116:LEU:HD22	2.22	0.69
1:A:122:ASN:HD21	1:A:307:ARG:HD2	1.58	0.69
1:A:55:VAL:HG12	1:A:117:PHE:HZ	1.58	0.68
1:A:104:LEU:O	1:A:108:LEU:HB2	1.93	0.67
2:B:53:ILE:N	2:B:53:ILE:HD13	2.09	0.67
2:B:51:VAL:HG12	2:B:53:ILE:CD1	2.25	0.67
1:A:73:LEU:HD22	1:A:74:THR:H	1.59	0.66
1:A:72:GLY:N	3:A:508:HOH:O	2.28	0.66
2:B:3:GLN:HG2	2:B:4:ILE:N	2.09	0.66
1:A:24:LYS:NZ	1:A:35:GLU:OE2	2.29	0.66
2:B:53:ILE:N	2:B:53:ILE:CD1	2.59	0.66
2:B:339:THR:O	2:B:343:GLN:HA	1.96	0.66
2:B:49:GLN:HB2	2:B:64:ARG:O	1.96	0.65
1:A:312:ARG:HG2	1:A:312:ARG:HH11	1.62	0.65
2:B:76:ALA:HB3	2:B:116:LEU:HD22	1.77	0.65

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:101:GLN:O	1:A:105:ILE:HG13	2.00	0.61
1:A:63:TYR:O	1:A:73:LEU:HD23	2.00	0.61
2:B:78:LYS:HD2	2:B:94:GLU:OE2	2.00	0.61
2:B:253:PRO:CD	2:B:256:ALA:HB3	2.29	0.61
1:A:312:ARG:HB3	3:A:414:HOH:O	1.99	0.61
1:A:11:LYS:HD2	1:A:166:LEU:HD23	1.83	0.61
1:A:11:LYS:HD2	1:A:166:LEU:CD2	2.31	0.60
1:A:122:ASN:ND2	1:A:307:ARG:HD2	2.16	0.60
2:B:253:PRO:HD2	2:B:256:ALA:CB	2.28	0.60
1:A:15:ASN:ND2	1:A:17:LYS:HG3	2.14	0.60
1:A:166:LEU:HD22	1:A:383:PHE:CG	2.37	0.60
2:B:345:PRO:HD2	3:B:495:HOH:O	2.02	0.59
1:A:15:ASN:OD1	1:A:162:ASN:HA	2.03	0.59
2:B:315:ILE:C	2:B:315:ILE:HD12	2.21	0.58
2:B:14:GLN:OE1	2:B:16:GLN:HG3	2.04	0.58
2:B:11:GLU:OE2	2:B:323:GLN:HG2	2.04	0.57
2:B:8:ARG:HH12	2:B:10:GLN:NE2	1.86	0.57
1:A:39:GLY:O	1:A:101:GLN:NE2	2.37	0.57
2:B:51:VAL:HG12	2:B:53:ILE:HD11	1.87	0.57
1:A:306:LEU:CD2	1:A:315:LEU:CD1	2.83	0.57
2:B:76:ALA:HB2	2:B:116:LEU:CD2	2.34	0.56
1:A:191:LEU:O	1:A:192:MET:HB2	2.05	0.56
1:A:98:THR:OG1	1:A:101:GLN:HG3	2.06	0.55
2:B:196:GLY:HA2	2:B:220:VAL:HB	1.87	0.55
1:A:294:LYS:CE	3:A:415:HOH:O	2.28	0.55
2:B:51:VAL:HG22	2:B:63:ARG:HG2	1.87	0.55
2:B:148:LEU:HD12	2:B:148:LEU:N	2.22	0.55
1:A:55:VAL:HG12	1:A:117:PHE:CZ	2.41	0.55
1:A:18:LEU:HD22	1:A:41:VAL:HG22	1.87	0.55
1:A:201:ALA:HA	1:A:217:ASN:O	2.08	0.54
2:B:335:ILE:H	2:B:335:ILE:CD1	2.09	0.54
2:B:276:ILE:HG12	2:B:282:ILE:HD12	1.90	0.54
1:A:383:PHE:O	1:A:384:ALA:C	2.47	0.53
1:A:370:ILE:HG13	2:B:66:ILE:HG22	1.91	0.53
1:A:18:LEU:HD22	1:A:41:VAL:CG2	2.39	0.52
1:A:312:ARG:HH11	1:A:312:ARG:CG	2.23	0.52
1:A:236:ARG:HG2	1:A:236:ARG:HH11	1.75	0.51
1:A:64:GLY:O	1:A:73:LEU:HD23	2.10	0.51
1:A:24:LYS:HZ2	1:A:35:GLU:CD	2.13	0.51
1:A:156:GLU:HG2	1:A:157:LYS:H	1.75	0.51
2:B:301:GLU:HG3	2:B:321:LYS:HD2	1.91	0.51
2:B:12:HIS:O	2:B:13:LEU:HB3	2.09	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:63:TYR:HB3	1:A:75:PHE:HB3	1.93	0.50
1:A:43:VAL:HG22	1:A:111:HIS:O	2.10	0.50
2:B:59:SER:O	2:B:61:PRO:HD3	2.12	0.50
2:B:255:GLU:CD	2:B:255:GLU:H	2.15	0.50
2:B:49:GLN:HA	2:B:66:ILE:HG12	1.95	0.49
2:B:59:SER:C	2:B:61:PRO:HD3	2.33	0.49
1:A:15:ASN:OD1	1:A:161:ARG:O	2.31	0.49
1:A:73:LEU:CD2	1:A:74:THR:H	2.25	0.49
1:A:85:GLN:OE1	1:A:88:PRO:HD2	2.13	0.48
1:A:181:GLN:CB	1:A:182:PRO:CD	2.92	0.48
1:A:241:ILE:HD13	1:A:241:ILE:N	2.28	0.48
1:A:8:VAL:HG11	1:A:104:LEU:HD11	1.95	0.48
1:A:147:LYS:HZ3	1:A:158:ILE:HD12	1.75	0.48
1:A:73:LEU:HD22	1:A:74:THR:N	2.28	0.48
1:A:182:PRO:O	1:A:202:SER:HB2	2.14	0.48
1:A:326:LYS:HE2	3:A:436:HOH:O	2.13	0.47
2:B:315:ILE:CD1	2:B:316:ILE:N	2.72	0.47
1:A:236:ARG:HG2	1:A:236:ARG:NH1	2.30	0.47
2:B:13:LEU:HD12	2:B:13:LEU:C	2.35	0.47
1:A:307:ARG:HB2	1:A:310:ALA:HB2	1.97	0.46
2:B:11:GLU:HG3	2:B:325:LEU:HD23	1.97	0.46
2:B:59:SER:HB2	3:B:369:HOH:O	2.15	0.46
1:A:191:LEU:HD12	1:A:191:LEU:HA	1.74	0.46
1:A:232:LYS:HB3	1:A:232:LYS:HE2	1.69	0.46
1:A:312:ARG:CG	1:A:312:ARG:NH1	2.80	0.45
2:B:301:GLU:CD	2:B:321:LYS:HE3	2.37	0.45
1:A:53:VAL:HG22	1:A:150:CYS:SG	2.57	0.45
2:B:8:ARG:NH1	2:B:10:GLN:CG	2.80	0.45
2:B:237:PRO:O	2:B:240:ASN:HB2	2.17	0.45
2:B:52:ILE:CD1	2:B:52:ILE:N	2.79	0.45
1:A:306:LEU:CD2	1:A:315:LEU:HD13	2.46	0.45
2:B:355:ASN:O	2:B:356:ASN:CB	2.57	0.45
2:B:52:ILE:N	2:B:52:ILE:HD13	2.32	0.44
2:B:50:VAL:HG22	2:B:52:ILE:HD12	2.00	0.44
1:A:7:ARG:HH21	1:A:9:PHE:HZ	1.66	0.44
1:A:24:LYS:NZ	1:A:35:GLU:CD	2.70	0.43
2:B:69:ASP:N	2:B:83:LYS:O	2.51	0.43
2:B:86:LYS:NZ	2:B:105:THR:O	2.39	0.43
1:A:181:GLN:HB3	1:A:182:PRO:CD	2.49	0.43
1:A:185:GLU:OE2	1:A:198:HIS:CE1	2.72	0.42
1:A:73:LEU:HD22	1:A:75:PHE:H	1.83	0.42
1:A:314:ILE:HD13	1:A:314:ILE:HG21	1.71	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:303:ILE:HD13	2:B:319:ASN:HB3	2.00	0.42
1:A:62:ARG:HG3	1:A:62:ARG:HH11	1.85	0.42
1:A:251:CYS:HA	1:A:252:PRO:HD3	1.90	0.42
1:A:6:THR:O	1:A:378:ILE:HA	2.20	0.42
1:A:173:TYR:CE1	1:A:347:LYS:HB3	2.55	0.42
1:A:306:LEU:HD23	1:A:315:LEU:CD1	2.50	0.42
2:B:30:LEU:C	2:B:30:LEU:HD23	2.40	0.42
1:A:15:ASN:HD21	1:A:17:LYS:CG	2.31	0.42
2:B:164:TRP:CZ3	2:B:185:SER:HB2	2.54	0.42
2:B:25:ILE:HG22	2:B:25:ILE:O	2.20	0.41
2:B:83:LYS:NZ	3:B:434:HOH:O	2.53	0.41
1:A:379:VAL:HG12	1:A:380:PHE:N	2.35	0.41
1:A:307:ARG:HE	1:A:307:ARG:HB3	1.62	0.41
2:B:53:ILE:H	2:B:53:ILE:HD13	1.83	0.41
1:A:56:THR:HG23	1:A:147:LYS:HB3	2.03	0.40
1:A:232:LYS:HG2	1:A:257:GLU:HB3	2.03	0.40
2:B:335:ILE:N	2:B:336:PRO:HD2	2.36	0.40
2:B:51:VAL:HG22	2:B:63:ARG:CG	2.52	0.40
2:B:3:GLN:HG2	2:B:4:ILE:H	1.85	0.40
1:A:156:GLU:HG2	1:A:157:LYS:N	2.36	0.40
1:A:193:SER:HB3	3:A:470:HOH:O	2.21	0.40
1:A:166:LEU:HD22	1:A:383:PHE:CD2	2.56	0.40
1:A:314:ILE:HG13	3:A:501:HOH:O	2.21	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	313/385 (81%)	292 (93%)	20 (6%)	1 (0%)	50	53
2	B	342/363 (94%)	329 (96%)	12 (4%)	1 (0%)	50	53
All	All	655/748 (88%)	621 (95%)	32 (5%)	2 (0%)	50	53

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	160	LYS
2	B	94	GLU

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	286/344 (83%)	280 (98%)	6 (2%)	66	78
2	B	300/310 (97%)	291 (97%)	9 (3%)	53	64
All	All	586/654 (90%)	571 (97%)	15 (3%)	59	70

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	73	LEU
1	A	86	SER
1	A	110	GLU
1	A	250	LYS
1	A	277	PHE
1	A	367	THR
2	B	14	GLN
2	B	43	LYS
2	B	52	ILE
2	B	53	ILE
2	B	67	SER
2	B	93	ILE
2	B	111	TRP
2	B	282	ILE
2	B	315	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA chains 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 carbohydrates 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 ⓘ

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	329/385 (85%)	0.81	65 (19%) 2 1	4, 22, 49, 69	9 (2%)
2	B	349/363 (96%)	0.35	18 (5%) 26 26	6, 21, 39, 56	0
All	All	678/748 (90%)	0.57	83 (12%) 5 4	4, 22, 45, 69	9 (1%)

All (83) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	379	VAL	7.0
1	A	150	CYS	5.9
1	A	43	VAL	5.9
1	A	269	CYS	5.8
1	A	8	VAL	5.5
2	B	341	VAL	5.4
1	A	279	ALA	4.9
1	A	73	LEU	4.8
2	B	176	ARG	4.8
1	A	152	GLU	4.6
2	B	238	THR	4.5
1	A	161	ARG	4.4
1	A	53	VAL	4.3
1	A	108	LEU	4.3
1	A	378	ILE	4.3
1	A	19	THR	4.2
2	B	4	ILE	4.2
1	A	162	ASN	4.1
2	B	177	VAL	4.1
1	A	380	PHE	4.0
1	A	104	LEU	4.0
1	A	106	LYS	3.7
1	A	111	HIS	3.7
1	A	87	PHE	3.6

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Mol	Chain	Res	Type	RSRZ
1	A	112	ALA	3.5
1	A	20	VAL	3.4
1	A	16	GLY	3.4
1	A	307	ARG	3.4
1	A	86	SER	3.2
1	A	367	THR	3.2
1	A	17	LYS	3.0
1	A	110	GLU	3.0
1	A	5	GLY	3.0
1	A	52	ARG	3.0
1	A	277	PHE	2.9
2	B	235	THR	2.9
1	A	40	VAL	2.9
1	A	105	ILE	2.8
2	B	335	ILE	2.8
1	A	84	VAL	2.8
2	B	43	LYS	2.8
1	A	14	PRO	2.8
1	A	371	GLU	2.8
1	A	6	THR	2.7
1	A	12	ALA	2.7
1	A	10	LYS	2.7
1	A	244	PHE	2.7
1	A	192	MET	2.7
1	A	308	GLU	2.7
2	B	339	THR	2.7
2	B	3	GLN	2.7
1	A	88	PRO	2.6
1	A	113	TYR	2.6
1	A	294	LYS	2.6
2	B	172	ALA	2.6
2	B	186	VAL	2.6
1	A	96	PRO	2.5
1	A	15	ASN	2.5
1	A	51	ARG	2.4
1	A	72	GLY	2.4
1	A	42	LEU	2.4
1	A	219	HIS	2.4
2	B	178	VAL	2.3
1	A	54	TYR	2.3
2	B	45	GLY	2.2
2	B	173	GLN	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	41	VAL	2.2
2	B	51	VAL	2.2
1	A	311	ASN	2.2
1	A	163	SER	2.2
1	A	253	VAL	2.2
1	A	323	VAL	2.2
2	B	272	VAL	2.2
1	A	99	ARG	2.1
1	A	21	TYR	2.1
1	A	155	GLU	2.1
1	A	164	VAL	2.1
1	A	89	PRO	2.0
1	A	95	LYS	2.0
1	A	372	LEU	2.0
1	A	216	VAL	2.0
1	A	158	ILE	2.0
2	B	154	ILE	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 ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.