Full wwPDB X-ray Structure Validation Report (i)

Sep 14, 2023 – 10:39 AM EDT

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 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 (i) 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 (1)) were used in the production of this report:

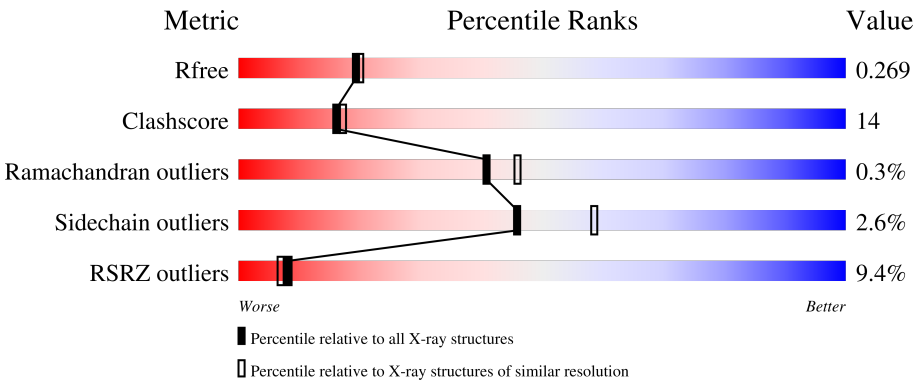
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

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}	130704	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (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 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 <=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	385	14% 62% 22% • 15%
2	B	363	3% 77% 16% • •

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 6221 atoms, of which 634 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 Beta-arrestin-1.

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

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

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

- Molecule 3 is water.

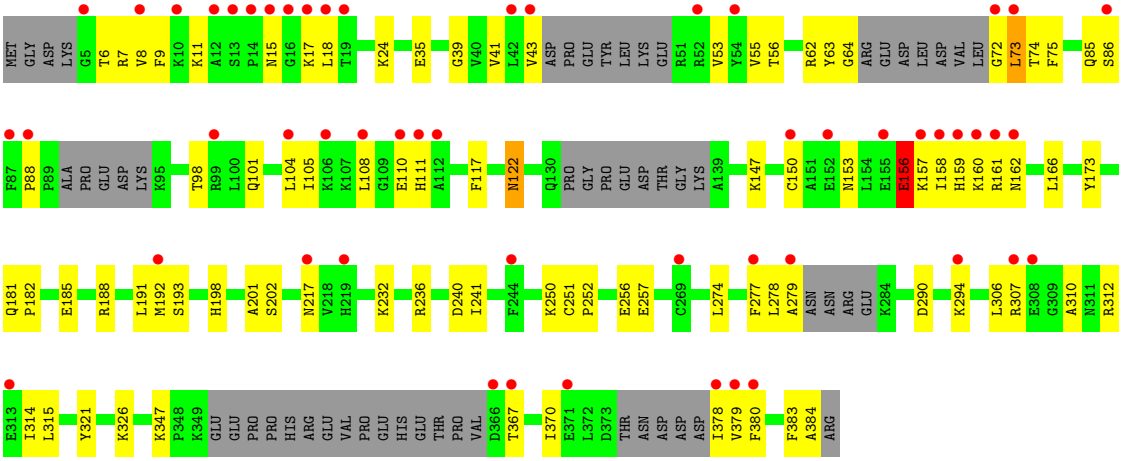
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	133	133	133	0	0
			Total	O		
3	B	143	143	143	0	0

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.

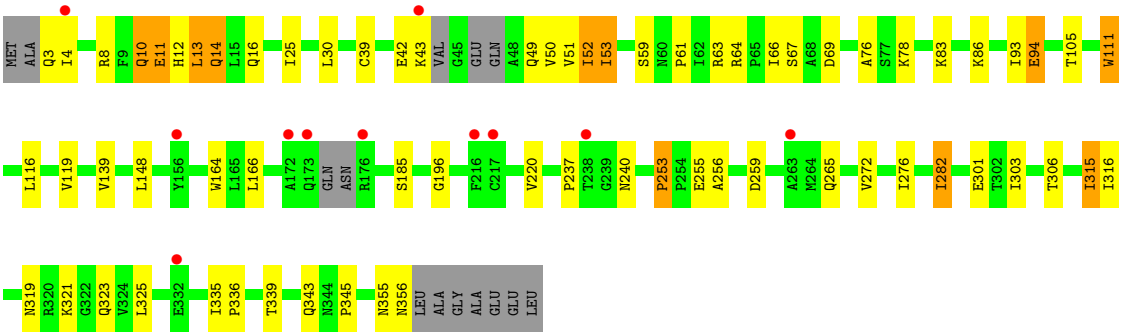
- Molecule 1: Beta-arrestin-1

Chain A:



- Molecule 2: Clathrin heavy chain 1

Chain B:



4 Data and refinement statistics (i)

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
$< I/\sigma(I) >$ [1]	2.89 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.4.0034	Depositor
R, R_{free}	0.200 , 0.251 0.232 , 0.269	Depositor DCC
R_{free} test set	3099 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å^2)	36.4	Xtriage
Anisotropy	0.044	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 49.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.011 for -h,l,k	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.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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	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(^{o})	Ideal(^{o})
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 Too-close contacts (i)

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	2582	322	2627	91	0
2	B	2729	312	2740	64	0
3	A	133	0	0	8	0
3	B	143	0	0	4	0
All	All	5587	634	5367	154	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 14.

All (154) 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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:147:LYS:NZ	1:A:158:ILE:CD1	2.06	1.18
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:HD12	2:B:335:ILE:H	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:312:ARG:HH11	1:A:312:ARG:HG2	1.62	0.65
2:B:76:ALA:HB3	2:B:116:LEU:HD22	1.77	0.65
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:11:LYS:HD2	1:A:166:LEU:HD23	1.83	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: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:HD12	2:B:315:ILE:C	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
1:A:306:LEU:CD2	1:A:315:LEU:CD1	2.83	0.57
2:B:51:VAL:HG12	2:B:53:ILE:HD11	1.87	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:N	2:B:148:LEU:HD12	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:276:ILE:HG12	2:B:282:ILE:HD12	1.90	0.54
2:B:335:ILE:H	2:B:335:ILE:CD1	2.09	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:24:LYS:HZ2	1:A:35:GLU:CD	2.13	0.51
1:A:64:GLY:O	1:A:73:LEU:HD23	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
1:A:43:VAL:HG22	1:A:111:HIS:O	2.10	0.50
1:A:63:TYR:HB3	1:A:75:PHE:HB3	1.93	0.50
2:B:59:SER:O	2:B:61:PRO:HD3	2.12	0.50
2:B:255:GLU:H	2:B:255:GLU:CD	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:N	1:A:241:ILE:HD13	2.28	0.48
1:A:8:VAL:HG11	1:A:104:LEU:HD11	1.95	0.48
1:A:73:LEU:HD22	1:A:74:THR:N	2.28	0.48
1:A:147:LYS:HZ3	1:A:158:ILE:HD12	1.75	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:C	2:B:13:LEU:HD12	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
1:A:53:VAL:HG22	1:A:150:CYS:SG	2.57	0.45
2:B:301:GLU:CD	2:B:321:LYS:HE3	2.37	0.45
2:B:8:ARG:NH1	2:B:10:GLN:CG	2.80	0.45
2:B:52:ILE:CD1	2:B:52:ILE:N	2.79	0.45
2:B:237:PRO:O	2:B:240:ASN:HB2	2.17	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
1:A:62:ARG:HH11	1:A:62:ARG:HG3	1.85	0.42
2:B:303:ILE:HD13	2:B:319:ASN:HB3	2.00	0.42
1:A:6:THR:O	1:A:378:ILE:HA	2.20	0.42
1:A:251:CYS:HA	1:A:252:PRO:HD3	1.90	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
1:A:379:VAL:HG12	1:A:380:PHE:N	2.35	0.41
2:B:83:LYS:NZ	3:B:434:HOH:O	2.53	0.41
1:A:307:ARG:HE	1:A:307:ARG:HB3	1.62	0.41
2:B:53:ILE:HD13	2:B:53:ILE:H	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
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
2:B:3:GLN:HG2	2:B:4:ILE:H	1.85	0.40
2:B:51:VAL:HG22	2:B:63:ARG:CG	2.52	0.40
2:B:335:ILE:N	2:B:336:PRO:HD2	2.36	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 (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	313/385 (81%)	292 (93%)	20 (6%)	1 (0%)	41	46
2	B	342/363 (94%)	329 (96%)	12 (4%)	1 (0%)	41	46
All	All	655/748 (88%)	621 (95%)	32 (5%)	2 (0%)	41	46

All (2) Ramachandran outliers are listed below:

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

5.3.2 Protein sidechains (i)

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%)	53	67
2	B	300/310 (97%)	291 (97%)	9 (3%)	41	53
All	All	586/654 (90%)	571 (97%)	15 (3%)	46	58

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

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Mol	Chain	Res	Type
2	B	315	ILE

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

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 monosaccharides 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 (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	329/385 (85%)	0.53	53 (16%) 1 1	4, 22, 49, 69	9 (2%)
2	B	349/363 (96%)	0.15	11 (3%) 47 45	6, 22, 39, 56	0
All	All	678/748 (90%)	0.33	64 (9%) 8 7	4, 22, 45, 69	9 (1%)

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	269	CYS	5.5
1	A	150	CYS	5.0
1	A	73	LEU	4.7
2	B	176	ARG	4.5
1	A	161	ARG	4.5
1	A	152	GLU	4.5
1	A	5	GLY	4.2
1	A	43	VAL	4.1
1	A	16	GLY	4.1
1	A	279	ALA	4.0
1	A	379	VAL	4.0
1	A	106	LYS	4.0
1	A	367	THR	3.9
1	A	162	ASN	3.8
1	A	111	HIS	3.6
1	A	108	LEU	3.6
1	A	380	PHE	3.5
1	A	192	MET	3.5
1	A	378	ILE	3.5
1	A	52	ARG	3.5
1	A	110	GLU	3.5
1	A	17	LYS	3.2
1	A	307	ARG	3.2
1	A	99	ARG	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	104	LEU	3.0
2	B	173	GLN	2.9
1	A	155	GLU	2.9
1	A	112	ALA	2.9
1	A	87	PHE	2.8
1	A	8	VAL	2.8
1	A	14	PRO	2.7
1	A	294	LYS	2.7
1	A	219	HIS	2.7
1	A	19	THR	2.6
1	A	42	LEU	2.6
1	A	72	GLY	2.6
1	A	371	GLU	2.6
1	A	15	ASN	2.5
1	A	13	SER	2.5
1	A	54	TYR	2.5
1	A	308	GLU	2.5
2	B	238	THR	2.4
2	B	332	GLU	2.3
1	A	160	LYS	2.3
1	A	12	ALA	2.3
1	A	18	LEU	2.3
2	B	217	CYS	2.3
2	B	43	LYS	2.3
1	A	158	ILE	2.3
1	A	366	ASP	2.2
1	A	244	PHE	2.2
1	A	313	GLU	2.2
1	A	10	LYS	2.2
1	A	86	SER	2.2
1	A	277	PHE	2.1
1	A	88	PRO	2.1
1	A	157	LYS	2.1
1	A	217	ASN	2.1
2	B	263	ALA	2.1
1	A	159	HIS	2.1
2	B	216	PHE	2.1
2	B	172	ALA	2.0
2	B	156	TYR	2.0
2	B	4	ILE	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 monosaccharides 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.