Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 10:39 AM GMT

PDB ID	:	3MKR
Title	:	Crystal structure of yeast alpha/epsilon-COP subcomplex of the COPI vesicular coat
Authors	:	Lee, C.; Goldberg, J.
Deposited on	:	2010-04-15
Resolution	:	2.60 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

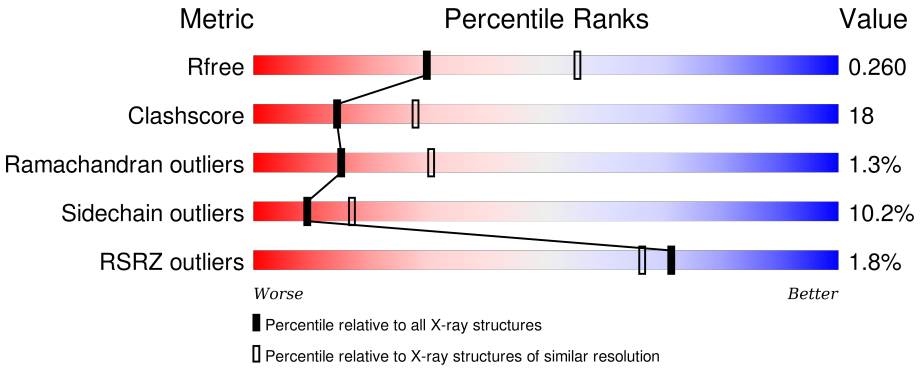
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)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

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.60 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	2328 (2.60-2.60)
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (2.60-2.60)

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 <=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	291	2% 60% 35% 5%
2	B	320	% 65% 30% 5%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4908 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 Coatomer subunit epsilon.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	291	Total	C	N	O	S	0	0	0
			2324	1460	403	449	12			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	94	HIS	ASP	VARIANT	UNP Q28104

- Molecule 2 is a protein called Coatomer subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	320	Total	C	N	O	S	0	0	0
			2550	1636	442	456	16			

- Molecule 3 is water.

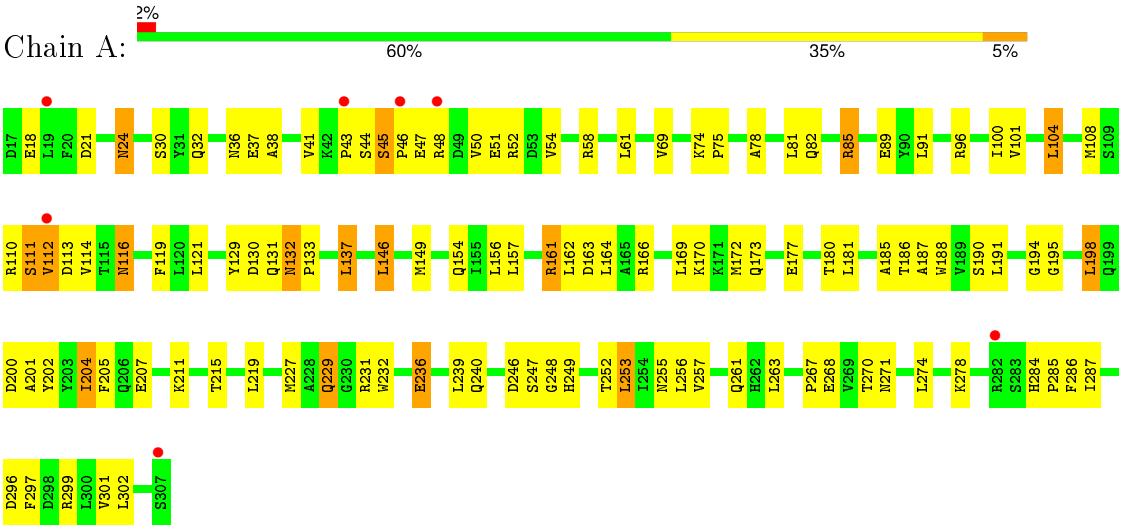
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	18	Total	O	0	0
			18	18		
3	B	16	Total	O	0	0
			16	16		

3 Residue-property plots (i)

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.

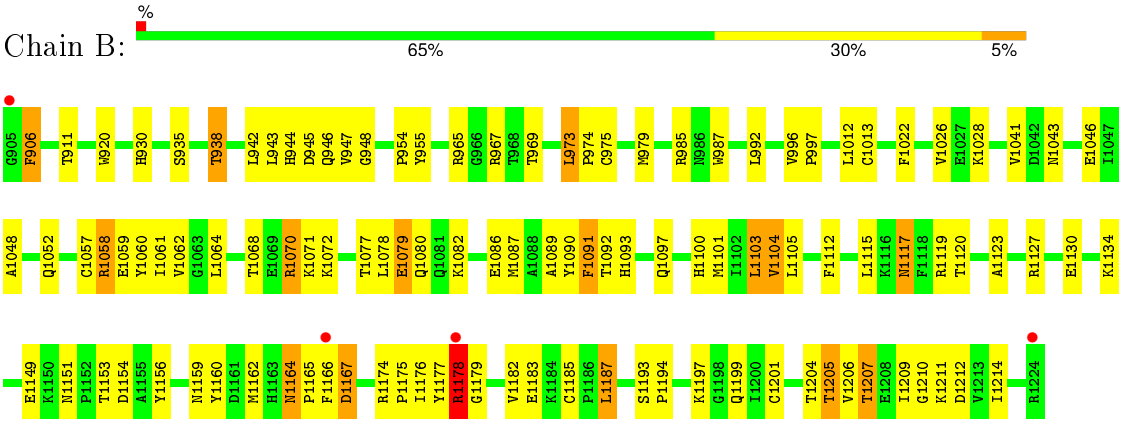
- Molecule 1: Coatomer subunit epsilon

Chain A:



- Molecule 2: Coatomer subunit alpha

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 4 2 2	Depositor
Cell constants a, b, c, α, β, γ	176.63Å 176.63Å 141.18Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.84 – 2.60 41.63 – 2.60	Depositor EDS
% Data completeness (in resolution range)	95.4 (29.84-2.60) 95.4 (41.63-2.60)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.04	Depositor
$< I/\sigma(I) >$ [1]	4.12 (at 2.61Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.222 , 0.263 0.221 , 0.260	Depositor DCC
R_{free} test set	1625 reflections (4.94%)	DCC
Wilson B-factor (Å^2)	51.0	Xtriage
Anisotropy	0.300	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 35.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Outliers	0 of 34363 reflections	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	4908	wwPDB-VP
Average B, all atoms (Å^2)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.60% 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, 0.375 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	0.38	0/2367	0.61	1/3201 (0.0%)
2	B	0.40	0/2607	0.68	0/3530
All	All	0.39	0/4974	0.64	1/6731 (0.0%)

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	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	229	GLN	N-CA-C	-5.58	95.94	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	202	TYR	Sidechain

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	2324	0	2287	87	0
2	B	2550	0	2608	99	0
3	A	18	0	0	0	0
3	B	16	0	0	0	0
All	All	4908	0	4895	177	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 18.

All (177) 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)
2:B:1185:CYS:SG	2:B:1207:THR:HG21	1.92	1.08
2:B:1178:ARG:HG2	2:B:1179:GLY:H	1.23	1.02
2:B:954:PRO:HG2	2:B:1206:VAL:HG12	1.54	0.87
2:B:1077:THR:HG22	2:B:1079:GLU:H	1.39	0.86
2:B:1204:THR:HG21	2:B:1207:THR:HB	1.57	0.85
1:A:74:LYS:HB2	1:A:75:PRO:HD2	1.59	0.84
1:A:255:ASN:HD21	2:B:969:THR:H	1.26	0.81
2:B:930:HIS:ND1	2:B:938:THR:HG21	1.98	0.78
2:B:1070:ARG:HH12	2:B:1071:LYS:HG3	1.49	0.78
1:A:38:ALA:O	1:A:52:ARG:HD2	1.84	0.77
2:B:954:PRO:HG3	2:B:1205:THR:HG23	1.67	0.77
2:B:1176:ILE:HG21	2:B:1182:VAL:HG13	1.67	0.77
2:B:1176:ILE:HG21	2:B:1182:VAL:CG1	2.17	0.75
1:A:170:LYS:HA	1:A:173:GLN:HB2	1.68	0.74
1:A:261:GLN:NE2	2:B:965:ARG:HH11	1.85	0.73
2:B:1178:ARG:HG2	2:B:1179:GLY:N	2.03	0.73
2:B:1057:CYS:O	2:B:1061:ILE:HG13	1.88	0.73
1:A:61:LEU:HD13	1:A:69:VAL:HG21	1.71	0.72
1:A:187:ALA:HB2	1:A:204:ILE:HG12	1.69	0.72
2:B:947:VAL:HG11	2:B:1187:LEU:HG	1.72	0.71
2:B:1164:ASN:HD22	2:B:1165:PRO:HD2	1.55	0.70
1:A:170:LYS:HA	1:A:173:GLN:HE21	1.57	0.70
2:B:1204:THR:CG2	2:B:1206:VAL:HG22	2.21	0.70
1:A:169:LEU:HD21	1:A:186:THR:HG22	1.72	0.70
2:B:947:VAL:HG12	2:B:948:GLY:N	2.07	0.69
2:B:930:HIS:ND1	2:B:938:THR:CG2	2.56	0.69
1:A:161:ARG:HD2	2:B:911:THR:O	1.92	0.68
2:B:1048:ALA:O	2:B:1052:GLN:HG3	1.94	0.68
2:B:954:PRO:HG2	2:B:1206:VAL:CG1	2.23	0.68
2:B:1089:ALA:O	2:B:1092:THR:HB	1.94	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:24:ASN:OD1	1:A:180:THR:HG21	1.94	0.68
2:B:1082:LYS:O	2:B:1086:GLU:HG3	1.95	0.67
1:A:85:ARG:HG3	1:A:85:ARG:HH11	1.59	0.67
2:B:1204:THR:HG22	2:B:1206:VAL:H	1.57	0.67
2:B:935:SER:OG	2:B:938:THR:HB	1.96	0.65
2:B:1022:PHE:CD2	2:B:1064:LEU:HD22	2.32	0.64
2:B:944:HIS:HD2	2:B:945:ASP:OD1	1.80	0.64
1:A:110:ARG:HG3	1:A:110:ARG:HH11	1.61	0.64
2:B:947:VAL:HG13	2:B:1187:LEU:HB2	1.79	0.64
1:A:154:GLN:OE1	2:B:973:LEU:HB2	1.98	0.64
1:A:188:TRP:HH2	2:B:973:LEU:HD13	1.63	0.64
2:B:1204:THR:O	2:B:1205:THR:HG22	1.98	0.64
2:B:996:VAL:HG22	2:B:997:PRO:HD2	1.81	0.63
1:A:261:GLN:HE22	2:B:965:ARG:HH11	1.46	0.63
2:B:1058:ARG:O	2:B:1062:VAL:HG13	1.99	0.62
2:B:1078:LEU:HG	2:B:1082:LYS:HE3	1.82	0.61
1:A:54:VAL:O	1:A:58:ARG:HG2	1.99	0.61
2:B:947:VAL:HG12	2:B:948:GLY:H	1.66	0.60
1:A:46:PRO:HG2	1:A:47:GLU:OE2	2.02	0.60
1:A:146:LEU:HD12	1:A:181:LEU:HD23	1.84	0.59
2:B:1177:TYR:HB3	2:B:1178:ARG:HD3	1.84	0.59
2:B:1204:THR:CG2	2:B:1207:THR:HB	2.32	0.59
1:A:188:TRP:CH2	2:B:973:LEU:HD13	2.38	0.59
2:B:954:PRO:CG	2:B:1206:VAL:HG12	2.31	0.58
2:B:1026:VAL:HG13	2:B:1061:ILE:HG23	1.86	0.58
1:A:110:ARG:HG3	1:A:110:ARG:NH1	2.19	0.58
1:A:112:VAL:O	1:A:112:VAL:HG12	2.04	0.57
2:B:1077:THR:HG22	2:B:1079:GLU:N	2.17	0.57
1:A:116:ASN:C	1:A:116:ASN:HD22	2.06	0.57
2:B:947:VAL:CG1	2:B:1187:LEU:HB2	2.34	0.57
2:B:1087:MET:HA	2:B:1090:TYR:CD2	2.40	0.56
2:B:1130:GLU:O	2:B:1130:GLU:HG2	2.04	0.56
1:A:89:GLU:OE2	1:A:96:ARG:HD3	2.06	0.56
2:B:1204:THR:HG22	2:B:1206:VAL:HG22	1.88	0.56
1:A:173:GLN:HG2	1:A:177:GLU:OE2	2.05	0.56
2:B:1086:GLU:OE2	2:B:1159:ASN:HB3	2.05	0.56
1:A:96:ARG:O	1:A:100:ILE:HG13	2.07	0.55
2:B:1201:CYS:CB	2:B:1207:THR:HG22	2.37	0.55
1:A:257:VAL:HG13	1:A:270:THR:HG23	1.88	0.55
1:A:161:ARG:HD3	1:A:164:LEU:HD12	1.89	0.55
1:A:149:MET:HB3	1:A:172:MET:HE2	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1078:LEU:O	2:B:1082:LYS:HG3	2.06	0.55
2:B:1058:ARG:HG3	2:B:1059:GLU:N	2.22	0.55
1:A:44:SER:O	1:A:45:SER:HB2	2.07	0.54
2:B:1093:HIS:HE1	2:B:1160:TYR:OH	1.90	0.54
1:A:255:ASN:HD21	2:B:969:THR:N	2.01	0.53
2:B:1060:TYR:OH	2:B:1097:GLN:HG2	2.06	0.53
2:B:985:ARG:HG2	2:B:996:VAL:HG13	1.90	0.53
2:B:1090:TYR:CZ	2:B:1175:PRO:HD3	2.44	0.53
2:B:1022:PHE:CE2	2:B:1064:LEU:HD22	2.44	0.53
2:B:1166:PHE:HE2	2:B:1177:TYR:CE2	2.27	0.53
1:A:78:ALA:O	1:A:82:GLN:HG2	2.09	0.53
2:B:1070:ARG:HH11	2:B:1070:ARG:HB3	1.75	0.52
2:B:1185:CYS:HG	2:B:1207:THR:HG21	1.74	0.52
1:A:24:ASN:HD21	1:A:180:THR:HG23	1.75	0.52
1:A:207:GLU:O	1:A:211:LYS:HG2	2.10	0.52
1:A:110:ARG:O	1:A:111:SER:HB2	2.09	0.52
2:B:1205:THR:HG22	2:B:1206:VAL:HG13	1.92	0.51
1:A:37:GLU:O	1:A:41:VAL:HG23	2.10	0.51
2:B:1077:THR:HG22	2:B:1078:LEU:N	2.25	0.51
2:B:1117:ASN:HA	2:B:1154:ASP:HA	1.93	0.51
1:A:85:ARG:HG3	1:A:85:ARG:NH1	2.25	0.51
2:B:1183:GLU:HG3	2:B:1197:LYS:HB2	1.93	0.51
2:B:1077:THR:HB	2:B:1080:GLN:H	1.76	0.51
1:A:61:LEU:CD1	1:A:69:VAL:HG21	2.39	0.51
2:B:1149:GLU:OE1	2:B:1149:GLU:HA	2.11	0.51
2:B:946:GLN:HA	2:B:1214:ILE:HB	1.94	0.50
1:A:161:ARG:NH1	1:A:164:LEU:HG	2.27	0.50
2:B:1105:LEU:HD11	2:B:1127:ARG:HD3	1.91	0.50
1:A:297:PHE:O	1:A:301:VAL:HG23	2.11	0.50
2:B:920:TRP:CG	2:B:997:PRO:HD3	2.47	0.50
2:B:1178:ARG:CG	2:B:1179:GLY:H	2.07	0.50
1:A:146:LEU:CD1	1:A:181:LEU:HB3	2.42	0.50
1:A:284:HIS:CG	1:A:285:PRO:HD2	2.47	0.50
1:A:161:ARG:NH1	1:A:163:ASP:OD2	2.45	0.50
1:A:32:GLN:HE21	1:A:36:ASN:ND2	2.10	0.50
2:B:1164:ASN:HD22	2:B:1165:PRO:CD	2.22	0.49
2:B:1178:ARG:HG3	2:B:1178:ARG:HH11	1.77	0.49
1:A:132:ASN:HD22	1:A:132:ASN:C	2.15	0.49
1:A:45:SER:OG	1:A:47:GLU:HG2	2.11	0.49
2:B:1197:LYS:HG3	2:B:1210:GLY:HA3	1.93	0.49
2:B:947:VAL:CG1	2:B:948:GLY:N	2.75	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:249:HIS:HD2	1:A:252:THR:OG1	1.95	0.49
1:A:204:ILE:HG13	1:A:205:PHE:N	2.28	0.49
1:A:24:ASN:HD21	1:A:180:THR:CG2	2.26	0.49
2:B:1112:PHE:CE2	2:B:1120:THR:HG21	2.48	0.49
2:B:1164:ASN:HD21	2:B:1178:ARG:HD2	1.77	0.49
1:A:169:LEU:HD11	1:A:186:THR:HG23	1.95	0.48
2:B:1201:CYS:HB2	2:B:1207:THR:HG22	1.94	0.48
1:A:185:ALA:HA	1:A:188:TRP:CE3	2.49	0.47
1:A:296:ASP:HB2	1:A:299:ARG:NH1	2.29	0.47
2:B:1013:CYS:SG	2:B:1028:LYS:HB3	2.54	0.47
1:A:190:SER:HB2	1:A:201:ALA:HB2	1.96	0.47
2:B:974:PRO:O	2:B:975:CYS:HB2	2.14	0.47
1:A:194:GLY:HA2	1:A:198:LEU:CD1	2.44	0.47
2:B:930:HIS:CE1	2:B:938:THR:HG21	2.50	0.47
1:A:30:SER:HB3	1:A:215:THR:HG22	1.96	0.46
1:A:113:ASP:OD1	1:A:114:VAL:N	2.49	0.46
2:B:1201:CYS:HB3	2:B:1207:THR:HG22	1.97	0.46
1:A:133:PRO:O	1:A:137:LEU:HB2	2.16	0.46
2:B:1043:ASN:OD1	2:B:1046:GLU:HB2	2.15	0.46
1:A:18:GLU:OE2	1:A:48:ARG:HG3	2.16	0.46
2:B:1151:ASN:OD1	2:B:1153:THR:HG22	2.17	0.45
2:B:1117:ASN:HD21	2:B:1156:TYR:H	1.63	0.45
2:B:1199:GLN:O	2:B:1209:ILE:HG13	2.17	0.45
1:A:236:GLU:HG2	1:A:263:LEU:HD11	1.98	0.45
2:B:1068:THR:O	2:B:1072:LYS:HG3	2.16	0.45
1:A:146:LEU:HD12	1:A:181:LEU:HB3	1.98	0.44
1:A:278:LYS:HG2	1:A:287:ILE:HD12	1.98	0.44
2:B:1123:ALA:HB1	2:B:1162:MET:HE1	1.98	0.44
1:A:274:LEU:HD12	1:A:274:LEU:HA	1.75	0.44
1:A:116:ASN:ND2	1:A:119:PHE:H	2.15	0.44
1:A:101:VAL:HG21	1:A:130:ASP:OD1	2.18	0.44
2:B:1064:LEU:O	2:B:1068:THR:HG23	2.17	0.43
2:B:1060:TYR:OH	2:B:1100:HIS:HD2	2.01	0.43
2:B:1115:LEU:O	2:B:1115:LEU:HG	2.18	0.43
1:A:284:HIS:CD2	1:A:286:PHE:H	2.37	0.43
2:B:1176:ILE:HD13	2:B:1182:VAL:HG12	2.01	0.43
1:A:200:ASP:O	1:A:204:ILE:HG23	2.19	0.43
2:B:1166:PHE:HB3	2:B:1167:ASP:H	1.42	0.43
1:A:267:PRO:HG2	1:A:268:GLU:OE2	2.18	0.43
2:B:1100:HIS:O	2:B:1104:VAL:HG13	2.18	0.43
1:A:46:PRO:O	1:A:50:VAL:HG23	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:161:ARG:HH11	1:A:161:ARG:HB3	1.84	0.43
1:A:227:MET:HB3	1:A:232:TRP:CZ3	2.55	0.42
1:A:21:ASP:HB3	1:A:37:GLU:OE1	2.19	0.42
2:B:1064:LEU:HD23	2:B:1091:PHE:HZ	1.85	0.42
1:A:110:ARG:O	1:A:111:SER:CB	2.67	0.42
1:A:129:TYR:C	1:A:131:GLN:H	2.22	0.42
1:A:246:ASP:OD1	1:A:247:SER:O	2.38	0.42
1:A:248:GLY:H	1:A:253:LEU:HD21	1.83	0.42
1:A:74:LYS:HB2	1:A:75:PRO:CD	2.41	0.42
1:A:229:GLN:OE1	1:A:231:ARG:NH1	2.48	0.42
1:A:146:LEU:CD1	1:A:181:LEU:HD23	2.50	0.41
2:B:947:VAL:CG1	2:B:948:GLY:H	2.32	0.41
1:A:104:LEU:O	1:A:108:MET:HB2	2.19	0.41
2:B:1103:LEU:HA	2:B:1103:LEU:HD12	1.92	0.41
1:A:161:ARG:HH11	1:A:161:ARG:CG	2.34	0.41
1:A:166:ARG:HG2	1:A:166:ARG:HH11	1.85	0.41
1:A:89:GLU:CD	1:A:96:ARG:HD3	2.41	0.41
1:A:195:GLY:O	1:A:198:LEU:HB2	2.21	0.41
1:A:229:GLN:HA	2:B:987:TRP:CE2	2.56	0.41
1:A:74:LYS:CB	1:A:75:PRO:HD2	2.33	0.40
2:B:1183:GLU:HG2	2:B:1194:PRO:HA	2.02	0.40
2:B:1119:ARG:HB3	2:B:1154:ASP:OD2	2.21	0.40
1:A:116:ASN:ND2	1:A:116:ASN:C	2.73	0.40
1:A:114:VAL:O	1:A:114:VAL:HG12	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	289/291 (99%)	269 (93%)	16 (6%)	4 (1%)	14 28

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
2	B	318/320 (99%)	285 (90%)	29 (9%)	4 (1%)	15 30
All	All	607/611 (99%)	554 (91%)	45 (7%)	8 (1%)	15 30

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	43	PRO
1	A	111	SER
2	B	1167	ASP
2	B	1178	ARG
1	A	45	SER
2	B	1211	LYS
2	B	906	PHE
1	A	112	VAL

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	249/249 (100%)	223 (90%)	26 (10%)	9 16
2	B	280/280 (100%)	252 (90%)	28 (10%)	9 18
All	All	529/529 (100%)	475 (90%)	54 (10%)	9 17

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	ASN
1	A	51	GLU
1	A	81	LEU
1	A	85	ARG
1	A	91	LEU
1	A	104	LEU
1	A	116	ASN
1	A	121	LEU

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Mol	Chain	Res	Type
1	A	132	ASN
1	A	137	LEU
1	A	146	LEU
1	A	156	LEU
1	A	157	LEU
1	A	161	ARG
1	A	162	LEU
1	A	191	LEU
1	A	198	LEU
1	A	204	ILE
1	A	219	LEU
1	A	236	GLU
1	A	239	LEU
1	A	240	GLN
1	A	253	LEU
1	A	256	LEU
1	A	271	ASN
1	A	302	LEU
2	B	906	PHE
2	B	938	THR
2	B	942	LEU
2	B	943	LEU
2	B	955	TYR
2	B	967	ARG
2	B	973	LEU
2	B	979	MET
2	B	992	LEU
2	B	1012	LEU
2	B	1041	VAL
2	B	1058	ARG
2	B	1070	ARG
2	B	1079	GLU
2	B	1091	PHE
2	B	1101	MET
2	B	1103	LEU
2	B	1104	VAL
2	B	1117	ASN
2	B	1134	LYS
2	B	1164	ASN
2	B	1174	ARG
2	B	1178	ARG
2	B	1187	LEU

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Mol	Chain	Res	Type
2	B	1193	SER
2	B	1205	THR
2	B	1207	THR
2	B	1212	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	ASN
1	A	39	GLN
1	A	116	ASN
1	A	132	ASN
1	A	173	GLN
1	A	249	HIS
1	A	255	ASN
1	A	261	GLN
1	A	271	ASN
2	B	923	ASN
2	B	944	HIS
2	B	946	GLN
2	B	951	GLN
2	B	1093	HIS
2	B	1100	HIS
2	B	1117	ASN
2	B	1139	GLN
2	B	1140	GLN
2	B	1163	HIS
2	B	1164	ASN

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 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, 95^{th} 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	291/291 (100%)	-0.06	7 (2%)	62	56	32, 48, 74, 90	0
2	B	320/320 (100%)	-0.08	4 (1%)	79	75	32, 51, 78, 90	0
All	All	611/611 (100%)	-0.07	11 (1%)	71	66	32, 50, 77, 90	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	112	VAL	5.8
2	B	1224	ARG	5.7
1	A	307	SER	4.3
1	A	19	LEU	3.2
2	B	1178	ARG	3.1
2	B	905	GLY	2.9
1	A	46	PRO	2.8
1	A	43	PRO	2.7
1	A	48	ARG	2.4
1	A	282	ARG	2.1
2	B	1166	PHE	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.