



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 20, 2023 – 12:41 PM EDT

PDB ID : 2HWW
Title : Structure of PIN domain of human SMG6
Authors : Glavan, F.; Behm-Ansmant, I.; Izaurralde, E.; Conti, E.
Deposited on : 2006-08-02
Resolution : 1.80 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 : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.35
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

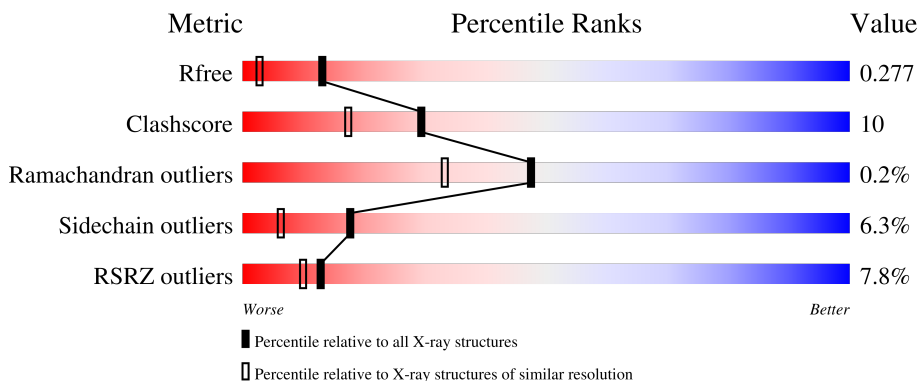
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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}	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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	181	
1	B	181	
1	C	181	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3943 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 Telomerase-binding protein EST1A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	160	1286	816	231	233	6	50	0	0
1	B	157	1272	809	228	230	5	45	0	0
1	C	144	1164	743	212	205	4	93	0	0

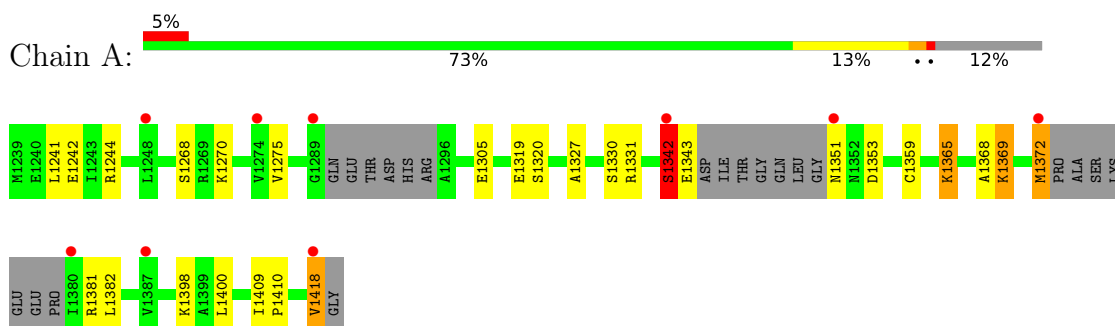
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	84	Total 84	O 84	0	0
2	B	82	Total 82	O 82	0	0
2	C	55	Total 55	O 55	0	0

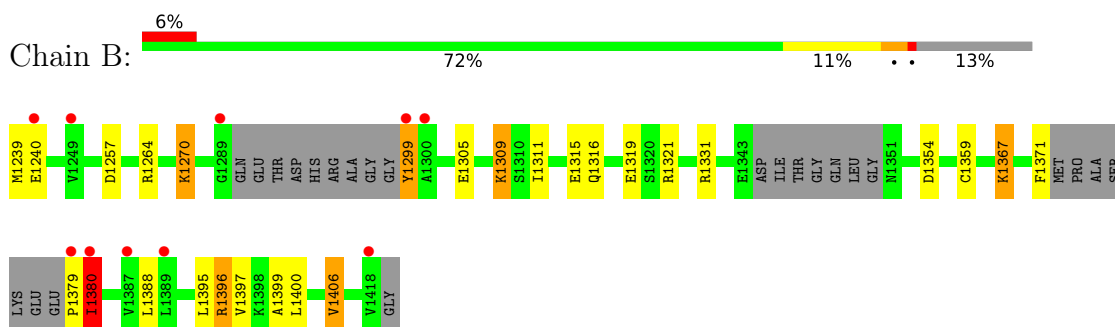
3 Residue-property plots [i](#)

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.

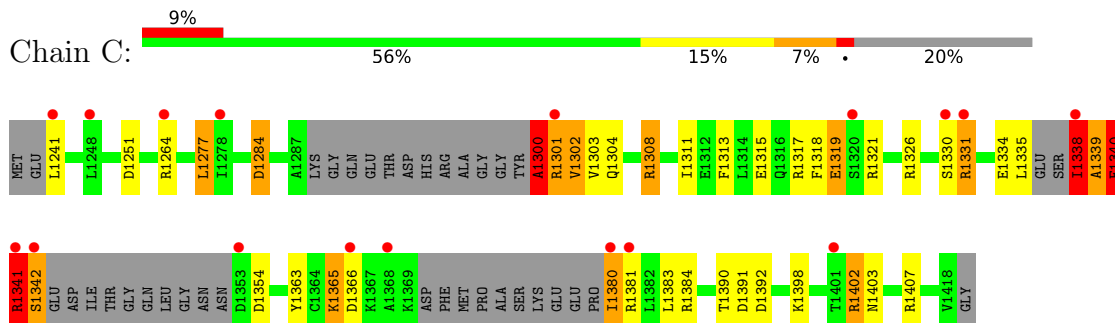
- Molecule 1: Telomerase-binding protein EST1A



- Molecule 1: Telomerase-binding protein EST1A



- Molecule 1: Telomerase-binding protein EST1A



4 Data and refinement statistics i

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	36.88Å 71.22Å 181.27Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.80 – 1.80 28.79 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.4 (28.80-1.80) 99.5 (28.79-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.72 (at 1.79Å)	Xtrriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.222 , 0.267 0.235 , 0.277	Depositor DCC
R_{free} test set	2290 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	24.6	Xtrriage
Anisotropy	0.085	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 52.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3943	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.49% 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	0.97	4/1301 (0.3%)	0.92	3/1750 (0.2%)
1	B	0.91	3/1288 (0.2%)	0.93	3/1734 (0.2%)
1	C	1.27	7/1176 (0.6%)	1.34	19/1583 (1.2%)
All	All	1.06	14/3765 (0.4%)	1.07	25/5067 (0.5%)

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
1	B	0	2
1	C	5	7
All	All	5	10

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	1331	ARG	C-N	-18.38	0.99	1.33
1	C	1330	SER	C-N	-15.15	0.99	1.34
1	C	1354	ASP	CB-CG	-10.87	1.28	1.51
1	A	1319	GLU	CB-CG	-9.23	1.34	1.52
1	B	1331	ARG	CB-CG	-8.67	1.29	1.52
1	B	1239	MET	C-N	-8.62	1.14	1.34
1	A	1381	ARG	CG-CD	-8.17	1.31	1.51
1	C	1321	ARG	CA-CB	-7.02	1.38	1.53
1	B	1406	VAL	CB-CG2	-6.28	1.39	1.52
1	C	1300	ALA	CA-CB	-6.13	1.39	1.52
1	C	1383	LEU	CB-CG	-5.94	1.35	1.52
1	C	1319	GLU	CG-CD	-5.33	1.44	1.51
1	A	1359	CYS	CB-SG	-5.07	1.73	1.81
1	A	1365	LYS	CB-CG	-5.03	1.39	1.52

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	1301	ARG	O-C-N	-16.62	96.10	122.70
1	C	1330	SER	O-C-N	-13.93	100.42	122.70
1	C	1331	ARG	O-C-N	-12.64	101.72	123.20
1	C	1330	SER	C-N-CA	11.73	151.02	121.70
1	C	1330	SER	CA-C-N	10.10	139.43	117.20
1	C	1331	ARG	CA-C-N	10.09	136.37	116.20
1	A	1381	ARG	CG-CD-NE	-8.30	94.37	111.80
1	C	1354	ASP	CB-CG-OD2	8.15	125.64	118.30
1	C	1341	ARG	N-CA-C	8.00	132.60	111.00
1	A	1369	LYS	CB-CG-CD	-7.22	92.82	111.60
1	C	1319	GLU	CB-CG-CD	7.02	133.16	114.20
1	C	1301	ARG	C-N-CA	6.87	138.88	121.70
1	C	1380	ILE	N-CA-C	6.62	128.86	111.00
1	B	1380	ILE	CG1-CB-CG2	6.56	125.83	111.40
1	C	1341	ARG	N-CA-CB	-6.19	99.45	110.60
1	B	1396	ARG	NE-CZ-NH2	6.13	123.37	120.30
1	C	1301	ARG	CA-C-N	5.94	130.26	117.20
1	C	1308	ARG	NE-CZ-NH1	5.92	123.26	120.30
1	A	1381	ARG	CB-CG-CD	5.88	126.88	111.60
1	B	1354	ASP	CB-CG-OD2	5.83	123.55	118.30
1	C	1300	ALA	C-N-CA	5.60	135.70	121.70
1	C	1339	ALA	N-CA-C	5.53	125.92	111.00
1	C	1366	ASP	CB-CG-OD1	-5.42	113.42	118.30
1	C	1284	ASP	CB-CG-OD2	5.28	123.05	118.30
1	C	1340	PHE	CB-CA-C	5.20	120.79	110.40

All (5) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	C	1338	ILE	CB
1	C	1339	ALA	CA
1	C	1340	PHE	CA
1	C	1380	ILE	CA
1	C	1390	THR	CB

All (10) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1342	SER	Peptide
1	B	1299	TYR	Peptide
1	B	1379	PRO	Peptide

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Mol	Chain	Res	Type	Group
1	C	1300	ALA	Peptide
1	C	1301	ARG	Mainchain,Peptide
1	C	1331	ARG	Mainchain
1	C	1338	ILE	Peptide
1	C	1340	PHE	Peptide
1	C	1402	ARG	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	1286	0	1328	18	1
1	B	1272	0	1315	15	4
1	C	1164	0	1221	36	1
2	A	84	0	0	0	0
2	B	82	0	0	5	4
2	C	55	0	0	3	1
All	All	3943	0	3864	68	6

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 10.

All (68) 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:1359:CYS:HB2	2:B:76:HOH:O	1.38	1.20
1:C:1390:THR:HG22	1:C:1392:ASP:H	1.31	0.95
1:C:1407:ARG:CZ	2:C:283:HOH:O	2.15	0.94
1:C:1339:ALA:O	1:C:1340:PHE:HB3	1.68	0.90
1:C:1300:ALA:HB2	1:C:1302:VAL:HG13	1.54	0.89
1:C:1300:ALA:CB	1:C:1302:VAL:HG13	2.14	0.77
1:A:1270:LYS:HD3	1:A:1418:VAL:HG21	1.66	0.77
1:C:1340:PHE:HD2	1:C:1342:SER:HG	1.32	0.77
1:A:1268:SER:HB2	1:A:1418:VAL:HG22	1.67	0.76
1:C:1251:ASP:OD2	1:C:1390:THR:HG23	1.86	0.75
1:A:1342:SER:HB2	1:A:1343:GLU:HB3	1.68	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1334:GLU:OE2	1:C:1363:TYR:OH	2.09	0.70
1:A:1268:SER:HB2	1:A:1418:VAL:CG2	2.22	0.70
1:C:1407:ARG:NE	2:C:283:HOH:O	2.21	0.69
1:C:1340:PHE:HD2	1:C:1342:SER:OG	1.75	0.68
1:A:1342:SER:CB	1:A:1343:GLU:HB3	2.26	0.65
1:B:1257:ASP:OD1	2:B:154:HOH:O	2.15	0.63
1:C:1241:LEU:HD21	1:C:1384:ARG:HH21	1.65	0.61
1:C:1340:PHE:CD2	1:C:1342:SER:OG	2.54	0.60
1:C:1390:THR:HG22	1:C:1392:ASP:N	2.09	0.60
1:B:1371:PHE:CD2	1:B:1380:ILE:HD13	2.37	0.58
1:B:1399:ALA:HB3	1:B:1406:VAL:HG21	1.86	0.57
1:A:1241:LEU:HD22	1:A:1382:LEU:HD23	1.86	0.56
1:B:1311:ILE:O	1:B:1315:GLU:HG3	2.04	0.56
1:A:1242:GLU:OE2	1:A:1244:ARG:NE	2.38	0.56
1:B:1270:LYS:HE3	2:B:289:HOH:O	2.04	0.56
1:A:1342:SER:HB2	1:A:1343:GLU:CB	2.37	0.54
1:A:1342:SER:HB2	1:A:1343:GLU:HA	1.90	0.54
1:C:1277:LEU:HD21	1:C:1342:SER:HB3	1.89	0.54
1:B:1371:PHE:HD2	1:B:1380:ILE:HD13	1.73	0.53
1:A:1241:LEU:CD2	1:A:1382:LEU:HD23	2.38	0.52
1:C:1304:GLN:O	1:C:1308:ARG:HG3	2.10	0.52
1:C:1318:PHE:HD2	1:C:1338:ILE:HG22	1.75	0.52
1:A:1368:ALA:HA	1:B:1397:VAL:HG21	1.92	0.52
1:C:1390:THR:HG22	1:C:1391:ASP:N	2.24	0.52
1:A:1353:ASP:OD1	1:A:1398:LYS:NZ	2.30	0.51
1:A:1369:LYS:O	1:A:1372:MET:HB3	2.11	0.50
1:A:1342:SER:HB2	1:A:1343:GLU:CA	2.41	0.50
1:C:1390:THR:CG2	1:C:1392:ASP:H	2.15	0.50
1:C:1391:ASP:C	1:C:1391:ASP:OD2	2.50	0.50
1:C:1300:ALA:HB1	1:C:1303:VAL:H	1.76	0.49
1:C:1335:LEU:HB3	1:C:1339:ALA:HB3	1.95	0.49
1:C:1338:ILE:HA	1:C:1339:ALA:HB3	1.95	0.49
1:C:1313:PHE:O	1:C:1317:ARG:HG2	2.12	0.48
1:C:1365:LYS:O	1:C:1365:LYS:CG	2.60	0.48
1:A:1268:SER:CB	1:A:1418:VAL:HG22	2.40	0.47
1:C:1341:ARG:O	1:C:1341:ARG:HG3	2.14	0.47
1:B:1367:LYS:HZ2	1:B:1367:LYS:HB3	1.82	0.45
1:C:1380:ILE:O	1:C:1381:ARG:HB2	2.18	0.44
1:B:1305:GLU:HG3	2:B:89:HOH:O	2.18	0.44
1:C:1338:ILE:CA	1:C:1339:ALA:CB	2.95	0.44
1:B:1305:GLU:OE1	1:B:1309:LYS:NZ	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1316:GLN:HA	1:B:1319:GLU:HG2	2.00	0.43
1:A:1275:VAL:O	1:A:1327:ALA:HA	2.19	0.43
1:C:1338:ILE:CA	1:C:1339:ALA:HB3	2.48	0.43
1:B:1399:ALA:CB	1:B:1406:VAL:CG2	2.98	0.42
1:B:1305:GLU:CG	2:B:89:HOH:O	2.68	0.42
1:C:1407:ARG:NH2	2:C:111:HOH:O	2.53	0.42
1:C:1315:GLU:O	1:C:1319:GLU:HG3	2.20	0.41
1:C:1390:THR:CG2	1:C:1391:ASP:N	2.84	0.41
1:A:1418:VAL:O	1:A:1418:VAL:CG1	2.69	0.41
1:B:1388:LEU:HD21	1:B:1395:LEU:HG	2.02	0.41
1:C:1251:ASP:OD2	1:C:1390:THR:CG2	2.63	0.41
1:A:1409:ILE:N	1:A:1410:PRO:CD	2.84	0.41
1:C:1311:ILE:HD13	1:C:1311:ILE:HA	1.74	0.41
1:C:1326:ARG:NH1	1:C:1334:GLU:OE2	2.54	0.41
1:C:1339:ALA:O	1:C:1340:PHE:CB	2.52	0.41
1:C:1264:ARG:HH11	1:C:1264:ARG:HD3	1.75	0.40

All (6) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1240:GLU:OE2	2:B:28:HOH:O[1_455]	0.53	1.67
1:B:1240:GLU:CD	2:B:28:HOH:O[1_455]	0.74	1.46
1:C:1398:LYS:NZ	2:C:215:HOH:O[1_655]	1.14	1.06
1:B:1240:GLU:OE1	2:B:28:HOH:O[1_455]	1.71	0.49
1:A:1305:GLU:OE2	1:A:1365:LYS:NZ[3_555]	1.95	0.25
1:B:1240:GLU:CG	2:B:28:HOH:O[1_455]	2.04	0.16

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	152/181 (84%)	148 (97%)	4 (3%)	0	100	100
1	B	149/181 (82%)	146 (98%)	3 (2%)	0	100	100
1	C	134/181 (74%)	127 (95%)	6 (4%)	1 (1%)	22	10
All	All	435/543 (80%)	421 (97%)	13 (3%)	1 (0%)	47	33

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	1403	ASN

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	142/159 (89%)	134 (94%)	8 (6%)	21	8
1	B	142/159 (89%)	133 (94%)	9 (6%)	18	6
1	C	130/159 (82%)	121 (93%)	9 (7%)	15	5
All	All	414/477 (87%)	388 (94%)	26 (6%)	18	6

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1320	SER
1	A	1330	SER
1	A	1331	ARG
1	A	1342	SER
1	A	1351	ASN
1	A	1372	MET
1	A	1400	LEU
1	A	1418	VAL
1	B	1264	ARG
1	B	1270	LYS
1	B	1299	TYR
1	B	1309	LYS

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Mol	Chain	Res	Type
1	B	1321	ARG
1	B	1367	LYS
1	B	1380	ILE
1	B	1396	ARG
1	B	1400	LEU
1	C	1277	LEU
1	C	1284	ASP
1	C	1302	VAL
1	C	1338	ILE
1	C	1340	PHE
1	C	1341	ARG
1	C	1342	SER
1	C	1365	LYS
1	C	1402	ARG

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	1281	ASN
1	C	1394	ASN

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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	C	2
1	B	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	1239:MET	C	1240:GLU	N	1.14
1	C	1331:ARG	C	1332:GLY	N	1.00
1	C	1330:SER	C	1331:ARG	N	0.99

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	160/181 (88%)	0.47	9 (5%) 24 19	15, 26, 38, 51	12 (7%)
1	B	157/181 (86%)	0.40	10 (6%) 19 15	15, 25, 42, 56	8 (5%)
1	C	144/181 (79%)	0.55	17 (11%) 4 3	20, 31, 46, 60	23 (15%)
All	All	461/543 (84%)	0.47	36 (7%) 13 10	15, 27, 42, 60	43 (9%)

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1372	MET	6.4
1	B	1289	GLY	5.0
1	C	1330	SER	4.1
1	C	1401	THR	4.0
1	B	1379	PRO	3.9
1	B	1418	VAL	3.6
1	A	1418	VAL	3.3
1	C	1381	ARG	3.2
1	B	1240	GLU	3.2
1	C	1248	LEU	3.1
1	B	1380	ILE	2.9
1	B	1300	ALA	2.8
1	A	1289	GLY	2.7
1	C	1353	ASP	2.7
1	A	1387	VAL	2.6
1	C	1264	ARG	2.6
1	C	1241	LEU	2.5
1	C	1368	ALA	2.4
1	B	1299	TYR	2.4
1	A	1380	ILE	2.4
1	A	1342	SER	2.3
1	C	1341	ARG	2.3
1	C	1380	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	1320	SER	2.2
1	C	1331	ARG	2.2
1	B	1387	VAL	2.2
1	A	1248	LEU	2.2
1	C	1366	ASP	2.2
1	C	1301	ARG	2.1
1	C	1338	ILE	2.1
1	C	1342	SER	2.1
1	B	1389	LEU	2.1
1	B	1249	VAL	2.1
1	A	1274	VAL	2.0
1	A	1351	ASN	2.0
1	C	1278	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.