



wwPDB X-ray Structure Validation Summary Report ⓘ

May 21, 2020 – 08:33 pm BST

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 wwPDB X-ray Structure Validation Summary 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 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 : 2.11
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) : 2.11

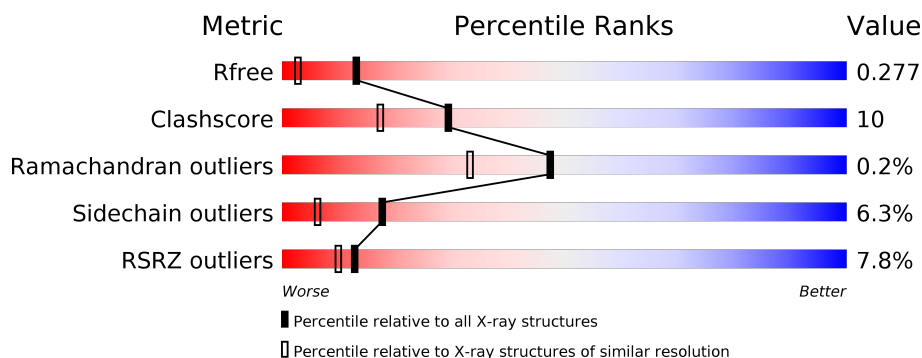
1 Overall quality at a glance

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 on 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	<div> <div>5%</div> <div> <div></div> <div>73%</div> <div>13%</div> <div>••</div> <div>12%</div> </div> </div>
1	B	181	<div> <div>6%</div> <div> <div></div> <div>72%</div> <div>11%</div> <div>••</div> <div>13%</div> </div> </div>
1	C	181	<div> <div>9%</div> <div> <div></div> <div>56%</div> <div>15%</div> <div>7%</div> <div>•</div> <div>20%</div> </div> </div>

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.

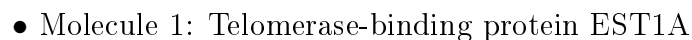
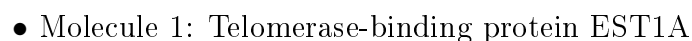
- Molecule 1 is a protein called Telomerase-binding protein EST1A.

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

- Molecule 2 is water.

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

- Molecule 1: Telomerase-binding protein EST1A



4 Data and refinement statistics

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)	Xtriage
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	Xtriage
Anisotropy	0.085	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 52.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
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

Xtriage'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

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	0.97	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

The worst 5 of 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

The worst 5 of 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

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

5 of 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
1	C	1300	ALA	Peptide
1	C	1301	ARG	Mainchain

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	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.

The worst 5 of 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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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

The worst 5 of 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

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

5 of 26 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	1309	LYS
1	B	1380	ILE
1	C	1365	LYS
1	B	1321	ARG
1	B	1367	LYS

Some 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 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](#)

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 [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	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, 47, 60	23 (15%)
All	All	461/543 (84%)	0.47	36 (7%) 13 10	15, 27, 42, 60	43 (9%)

The worst 5 of 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

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 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.