



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 07:22 AM UTC

PDB ID : 5EJC / pdb_00005ejc
Title : Crystal structural of the TSC1-TBC1D7 complex
Authors : Wang, Z.; Qin, J.; Gong, W.; Xu, W.
Deposited on : 2015-11-01
Resolution : 3.10 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 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:

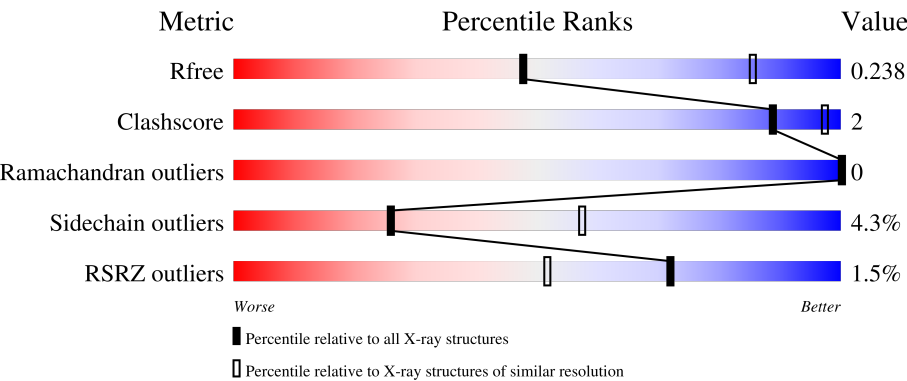
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

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 3.10 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}	180053	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)

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	276	% 92%5%.
1	B	276	% 91%5%.
2	C	55	44%13%44%
2	D	55	2%51%11%38%
2	E	55	4%47%7%.42%

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Mol	Chain	Length	Quality of chain
2	F	55	 A horizontal bar chart showing the quality of chain F. The bar is divided into three segments: a green segment representing 58%, a yellow segment representing 11%, and a grey segment representing 29%. A small black dot is located on the yellow segment.

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5467 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 TBC1 domain family member 7.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	268	Total	C	N	O	S	Se	0	0	0
			2179	1410	365	389	7	8			
1	B	267	Total	C	N	O	S	Se	0	0	0
			2169	1403	364	387	7	8			

- Molecule 2 is a protein called Hamartin.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	31	Total	C	N	O	0	0	0
			256	162	45	49			
2	D	34	Total	C	N	O	0	0	0
			278	175	49	54			
2	E	32	Total	C	N	O	0	0	0
			265	167	46	52			
2	F	39	Total	C	N	O	0	0	0
			320	203	56	61			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	938	GLY	-	expression tag	UNP Q92574
D	938	GLY	-	expression tag	UNP Q92574
E	938	GLY	-	expression tag	UNP Q92574
F	938	GLY	-	expression tag	UNP Q92574

● Molecule 1: TBC1 domain family member 7



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|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|------|------|------|------|------|------|------|------|------|-----|-----|-----|-----|-----|-----|
| GLY | PHE | ARG | G21 | G22 | E23 | L67 | H71 | H74 | M78 | R81 | K82 | V88 | F97 | T102 | L114 | R121 | D150 | R156 | M194 | A198 | P199 | G287 | THR | PRO | VAL | HIS | SER | SER |
|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|------|------|------|------|------|------|------|------|------|-----|-----|-----|-----|-----|-----|

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|-----|------|------|------|------|------|------|------|------|------|------|------|------|
| GLY | G939 | S946 | R947 | Y948 | K952 | R953 | Y954 | T955 | Q956 | Y957 | L960 | L969 | GLU | LYS | ASP | GLY | LEU | LEU | LYS | LYS | LEU | GLU | GLU | GLU | LYS | LYS | ALA | GLU | ALA | ALA | ALA | ALA | ALA | GLU | GLU | ABC |
|-----|------|------|------|------|------|------|------|------|------|------|------|------|

- [illegible]

- GLY
 G939
 G940
 S946
 R947
 R953
 V957
 L960
 E961
 E970
 LYS
 ASP
 GLY
 LEU
 LEU
 LEU
 LYS
 LYS
 LEU
 LEU
 GLU
 GLU
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 LYS
 ALA
 ALA
 GLU
 ALA
 ALA
 ALA
 ALA
 ALA
 GLU
 GLU
 ARG

- 
- WORLD WIDE
PDB
PROTEIN DATA BANK

Chain F: 58% 11% • 29%

GLY	S939	S946	R947	Y948	R953	V957	L960	L969	L975	R976	R977	ILEJ	GLU	GLU	GLU	GLU	LYS	ALA	GLU	ALA	ALA	GLU	ALA	ALA	ALA	GLU	GLU	ARG
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4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	183.40Å 66.70Å 98.37Å 90.00° 91.91° 90.00°	Depositor
Resolution (Å)	50.00 – 3.10 50.00 – 3.10	Depositor EDS
% Data completeness (in resolution range)	96.7 (50.00-3.10) 96.7 (50.00-3.10)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.01 (at 3.06Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.220 , 0.239 0.220 , 0.238	Depositor DCC
R_{free} test set	1101 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	82.8	Xtriage
Anisotropy	0.620	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 41.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.026 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5467	wwPDB-VP
Average B, all atoms (Å ²)	100.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.20% 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	0.42	0/2224	0.73	1/2998 (0.0%)
1	B	0.41	0/2213	0.74	1/2981 (0.0%)
2	C	0.58	0/258	0.84	0/345
2	D	0.55	0/280	1.00	0/373
2	E	0.53	0/267	0.93	1/357 (0.3%)
2	F	0.54	0/322	0.90	1/428 (0.2%)
All	All	0.44	0/5564	0.78	4/7482 (0.1%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	947	ARG	CG-CD-NE	6.45	126.19	112.00
1	A	194	MSE	CB-CA-C	-5.47	101.71	110.79
1	B	194	MSE	CB-CA-C	5.22	120.90	109.99
2	F	960	LEU	N-CA-CB	5.03	118.08	110.28

There are no chirality outliers.

There are no planarity outliers.

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	2179	0	2221	6	0
1	B	2169	0	2210	8	0
2	C	256	0	257	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	278	0	279	5	0
2	E	265	0	263	5	0
2	F	320	0	331	3	0
All	All	5467	0	5561	21	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 2.

The worst 5 of 21 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:114:LEU:HD12	2:D:957:VAL:HG11	1.78	0.64
2:E:947:ARG:NH2	2:F:948:TYR:CD2	2.73	0.57
1:A:71:HIS:HA	1:A:74:HIS:CD2	2.42	0.55
1:B:71:HIS:HA	1:B:74:HIS:CD2	2.41	0.55
1:B:121:ARG:NE	1:B:121:ARG:O	2.40	0.55

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	266/276 (96%)	256 (96%)	10 (4%)	0	100	100
1	B	265/276 (96%)	255 (96%)	10 (4%)	0	100	100
2	C	29/55 (53%)	29 (100%)	0	0	100	100
2	D	32/55 (58%)	32 (100%)	0	0	100	100
2	E	30/55 (54%)	30 (100%)	0	0	100	100
2	F	37/55 (67%)	37 (100%)	0	0	100	100
All	All	659/772 (85%)	639 (97%)	20 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	245/243 (101%)	236 (96%)	9 (4%)	30	61
1	B	243/243 (100%)	236 (97%)	7 (3%)	37	66
2	C	26/43 (60%)	22 (85%)	4 (15%)	2	12
2	D	28/43 (65%)	27 (96%)	1 (4%)	31	62
2	E	27/43 (63%)	25 (93%)	2 (7%)	13	40
2	F	33/43 (77%)	30 (91%)	3 (9%)	9	32
All	All	602/658 (92%)	576 (96%)	26 (4%)	26	57

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

Mol	Chain	Res	Type
1	B	121	ARG
2	C	952	LYS
2	F	969	LEU
2	C	946	SER
2	C	955	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 8 such sidechains are listed below:

Mol	Chain	Res	Type
2	F	940	GLN
2	E	940	GLN
1	B	182	ASN
1	B	47	GLN
1	B	283	HIS

5.3.3 RNA ⓘ

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 oligosaccharides 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	260/276 (94%)	-0.14	3 (1%) 76 58	68, 96, 125, 183	0
1	B	259/276 (93%)	-0.16	4 (1%) 72 52	68, 96, 129, 147	0
2	C	31/55 (56%)	-0.00	0 100 100	85, 102, 129, 144	0
2	D	34/55 (61%)	0.01	1 (2%) 53 32	76, 100, 152, 162	0
2	E	32/55 (58%)	0.03	2 (6%) 26 14	74, 95, 130, 134	0
2	F	39/55 (70%)	0.40	0 100 100	78, 101, 155, 162	0
All	All	655/772 (84%)	-0.09	10 (1%) 72 52	68, 96, 132, 183	0

The worst 5 of 10 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	289	PRO	4.6
1	A	288	THR	3.2
2	E	939	GLY	3.0
2	D	938	GLY	2.8
1	A	22	VAL	2.5

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 oligosaccharides 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.