



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 01:39 PM UTC

PDB ID : 7X39 / pdb_00007x39
Title : Structure of CIZ1 bound ERH
Authors : Wang, X.; Xu, C.
Deposited on : 2022-02-28
Resolution : 2.85 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	:	4-5-2 with Phenix2.0
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

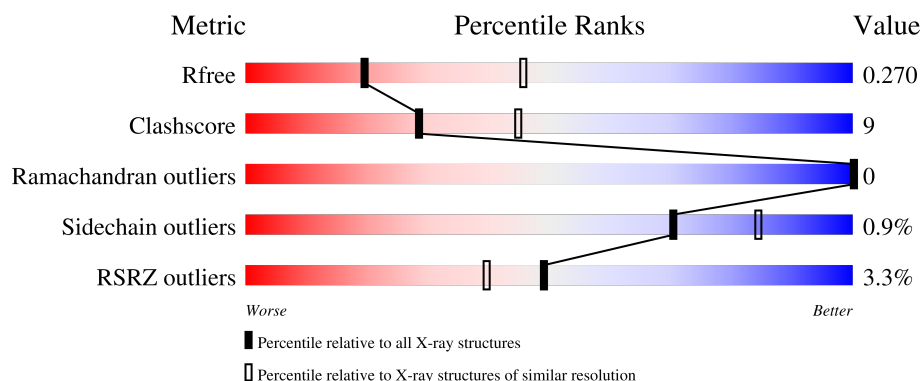
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 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	1407 (2.88-2.84)
Clashscore	190562	1446 (2.88-2.84)
Ramachandran outliers	187476	1406 (2.88-2.84)
Sidechain outliers	187428	1407 (2.88-2.84)
RSRZ outliers	180081	1408 (2.88-2.84)

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	158	7% 64% 16% 19%
1	B	158	58% 16% 26%
1	C	158	7% 55% 19% 26%
1	D	158	3% 54% 20% 26%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 3828 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 Enhancer of rudimentary homolog,Cip1-interacting zinc finger protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	128	Total	C	N	O	S	0	0	0
			1012	640	167	198	7			
1	B	117	Total	C	N	O	S	0	0	0
			943	595	154	187	7			
1	C	117	Total	C	N	O	S	0	0	0
			933	591	151	184	7			
1	D	117	Total	C	N	O	S	0	0	0
			940	594	154	185	7			

There are 92 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP P84090
A	0	HIS	-	expression tag	UNP P84090
A	105	GLY	-	linker	UNP P84090
A	106	SER	-	linker	UNP P84090
A	107	SER	-	linker	UNP P84090
A	108	GLY	-	linker	UNP P84090
A	109	SER	-	linker	UNP P84090
A	110	SER	-	linker	UNP P84090
A	111	GLY	-	linker	UNP P84090
A	112	SER	-	linker	UNP P84090
A	113	SER	-	linker	UNP P84090
A	114	GLY	-	linker	UNP P84090
A	115	SER	-	linker	UNP P84090
A	116	SER	-	linker	UNP P84090
A	117	GLY	-	linker	UNP P84090
A	118	SER	-	linker	UNP P84090
A	119	SER	-	linker	UNP P84090
A	120	GLY	-	linker	UNP P84090
A	121	SER	-	linker	UNP P84090
A	122	SER	-	linker	UNP P84090

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	123	GLY	-	linker	UNP P84090
A	124	SER	-	linker	UNP P84090
A	125	SER	-	linker	UNP P84090
B	-1	GLY	-	expression tag	UNP P84090
B	0	HIS	-	expression tag	UNP P84090
B	105	GLY	-	linker	UNP P84090
B	106	SER	-	linker	UNP P84090
B	107	SER	-	linker	UNP P84090
B	108	GLY	-	linker	UNP P84090
B	109	SER	-	linker	UNP P84090
B	110	SER	-	linker	UNP P84090
B	111	GLY	-	linker	UNP P84090
B	112	SER	-	linker	UNP P84090
B	113	SER	-	linker	UNP P84090
B	114	GLY	-	linker	UNP P84090
B	115	SER	-	linker	UNP P84090
B	116	SER	-	linker	UNP P84090
B	117	GLY	-	linker	UNP P84090
B	118	SER	-	linker	UNP P84090
B	119	SER	-	linker	UNP P84090
B	120	GLY	-	linker	UNP P84090
B	121	SER	-	linker	UNP P84090
B	122	SER	-	linker	UNP P84090
B	123	GLY	-	linker	UNP P84090
B	124	SER	-	linker	UNP P84090
B	125	SER	-	linker	UNP P84090
C	-1	GLY	-	expression tag	UNP P84090
C	0	HIS	-	expression tag	UNP P84090
C	105	GLY	-	linker	UNP P84090
C	106	SER	-	linker	UNP P84090
C	107	SER	-	linker	UNP P84090
C	108	GLY	-	linker	UNP P84090
C	109	SER	-	linker	UNP P84090
C	110	SER	-	linker	UNP P84090
C	111	GLY	-	linker	UNP P84090
C	112	SER	-	linker	UNP P84090
C	113	SER	-	linker	UNP P84090
C	114	GLY	-	linker	UNP P84090
C	115	SER	-	linker	UNP P84090
C	116	SER	-	linker	UNP P84090
C	117	GLY	-	linker	UNP P84090
C	118	SER	-	linker	UNP P84090

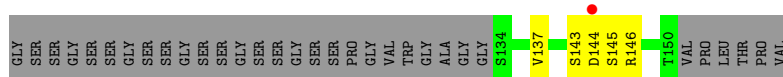
Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	119	SER	-	linker	UNP P84090
C	120	GLY	-	linker	UNP P84090
C	121	SER	-	linker	UNP P84090
C	122	SER	-	linker	UNP P84090
C	123	GLY	-	linker	UNP P84090
C	124	SER	-	linker	UNP P84090
C	125	SER	-	linker	UNP P84090
D	-1	GLY	-	expression tag	UNP P84090
D	0	HIS	-	expression tag	UNP P84090
D	105	GLY	-	linker	UNP P84090
D	106	SER	-	linker	UNP P84090
D	107	SER	-	linker	UNP P84090
D	108	GLY	-	linker	UNP P84090
D	109	SER	-	linker	UNP P84090
D	110	SER	-	linker	UNP P84090
D	111	GLY	-	linker	UNP P84090
D	112	SER	-	linker	UNP P84090
D	113	SER	-	linker	UNP P84090
D	114	GLY	-	linker	UNP P84090
D	115	SER	-	linker	UNP P84090
D	116	SER	-	linker	UNP P84090
D	117	GLY	-	linker	UNP P84090
D	118	SER	-	linker	UNP P84090
D	119	SER	-	linker	UNP P84090
D	120	GLY	-	linker	UNP P84090
D	121	SER	-	linker	UNP P84090
D	122	SER	-	linker	UNP P84090
D	123	GLY	-	linker	UNP P84090
D	124	SER	-	linker	UNP P84090
D	125	SER	-	linker	UNP P84090

- Molecule 1: Enhancer of rudimentary homolog, Cip1-interacting zinc finger protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	44.45Å 48.67Å 73.88Å 85.83° 75.08° 63.31°	Depositor
Resolution (Å)	38.08 – 2.85 38.08 – 2.85	Depositor EDS
% Data completeness (in resolution range)	96.5 (38.08-2.85) 96.5 (38.08-2.85)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.80 (at 2.85Å)	Xtriage
Refinement program	PHENIX 1.14 _3260	Depositor
R, R_{free}	0.218 , 0.267 0.218 , 0.270	Depositor DCC
R_{free} test set	695 reflections (5.58%)	wwPDB-VP
Wilson B-factor (Å ²)	40.6	Xtriage
Anisotropy	0.857	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 36.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.055 for h,h-k,h-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	3828	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.27% 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	0.33	0/1033	0.52	2/1402 (0.1%)
1	B	0.15	0/961	0.33	0/1302
1	C	0.29	0/951	0.43	0/1290
1	D	0.24	0/958	0.43	1/1298 (0.1%)
All	All	0.26	0/3903	0.43	3/5292 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	151	VAL	CA-C-N	6.66	126.44	119.05
1	A	151	VAL	C-N-CA	6.66	126.44	119.05
1	D	144	ASP	N-CA-C	-6.15	105.06	113.30

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	1012	0	979	16	0
1	B	943	0	911	14	0
1	C	933	0	895	18	0
1	D	940	0	909	20	0
All	All	3828	0	3694	66	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 9.

All (66) 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:D:58:PHE:HZ	1:D:92:TYR:HB2	1.65	0.60
1:C:95:LEU:HD22	1:C:148:PHE:HZ	1.67	0.59
1:A:51:THR:HG22	1:A:138:THR:HG23	1.84	0.59
1:B:70:LEU:HB3	1:B:79:TYR:HB3	1.86	0.58
1:A:80:GLN:HG2	1:A:82:TYR:CE2	2.40	0.57
1:A:134:SER:HB2	1:A:154:THR:HG23	1.85	0.57
1:D:73:ARG:HD2	1:D:80:GLN:HG3	1.88	0.55
1:C:20:ALA:HA	1:D:17:ARG:NH2	2.23	0.54
1:D:145:SER:OG	1:D:146:ARG:N	2.40	0.54
1:A:134:SER:HB2	1:A:154:THR:CG2	2.39	0.53
1:B:9:GLN:HB3	1:B:66:ASP:HB3	1.90	0.53
1:A:70:LEU:HB3	1:A:79:TYR:HB3	1.90	0.53
1:A:64:LEU:O	1:A:84:LYS:NZ	2.43	0.52
1:A:18:THR:OG1	1:A:19:TYR:N	2.43	0.52
1:C:33:CYS:HB2	1:C:146:ARG:HH21	1.74	0.52
1:A:54:ILE:HD13	1:A:95:LEU:HD12	1.93	0.50
1:C:18:THR:HA	1:D:18:THR:HA	1.94	0.50
1:A:139:ILE:HD12	1:A:145:SER:O	2.12	0.50
1:C:54:ILE:HG12	1:C:92:TYR:CE1	2.46	0.50
1:A:51:THR:HA	1:A:137:VAL:O	2.13	0.49
1:C:69:CYS:SG	1:C:82:TYR:HB2	2.53	0.49
1:C:61:ILE:HG23	1:C:67:LEU:HD23	1.95	0.49
1:B:138:THR:HG22	1:B:147:ALA:HB3	1.92	0.49
1:C:41:LYS:HD3	1:C:141:GLN:HG2	1.93	0.48
1:C:57:LEU:HD21	1:C:91:ILE:HG21	1.95	0.48
1:B:58:PHE:HB3	1:B:88:LYS:HE3	1.96	0.48
1:D:137:VAL:HG11	1:D:146:ARG:HD3	1.95	0.48
1:D:73:ARG:HD2	1:D:80:GLN:CG	2.44	0.48
1:D:27:GLU:HA	1:D:30:GLU:HB2	1.95	0.47
1:B:61:ILE:HA	1:B:64:LEU:HD12	1.96	0.47
1:C:73:ARG:HB3	1:C:75:ASP:OD1	2.15	0.47
1:D:69:CYS:HB2	1:D:87:ILE:HD11	1.97	0.46
1:C:92:TYR:CE2	1:C:96:ARG:HD2	2.50	0.46
1:D:70:LEU:HA	1:D:80:GLN:O	2.15	0.46
1:B:51:THR:HA	1:B:138:THR:HA	1.97	0.45
1:D:29:MET:HE1	1:D:87:ILE:HG12	1.97	0.45
1:C:14:PRO:O	1:C:17:ARG:HG3	2.16	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:25:VAL:O	1:D:29:MET:HG2	2.18	0.44
1:A:44:ASN:O	1:A:141:GLN:NE2	2.50	0.44
1:B:50:ILE:O	1:B:139:ILE:N	2.36	0.44
1:B:73:ARG:O	1:B:77:GLN:N	2.50	0.44
1:C:4:THR:HG23	1:C:71:VAL:HG22	2.00	0.43
1:D:38:GLU:HA	1:D:41:LYS:HE2	2.01	0.43
1:C:80:GLN:HG2	1:C:82:TYR:CZ	2.54	0.43
1:B:86:TRP:CH2	1:B:90:LYS:HG3	2.54	0.43
1:B:36:TYR:CZ	1:B:40:LEU:HD11	2.54	0.43
1:A:49:SER:HB3	1:A:140:LEU:HD23	2.00	0.42
1:B:49:SER:HB2	1:B:140:LEU:HD23	2.01	0.42
1:D:36:TYR:CD2	1:D:57:LEU:HD12	2.54	0.42
1:A:92:TYR:OH	1:A:96:ARG:NH1	2.52	0.42
1:D:73:ARG:HB3	1:D:75:ASP:OD1	2.19	0.42
1:A:139:ILE:HG13	1:A:140:LEU:N	2.35	0.42
1:B:73:ARG:HD2	1:B:80:GLN:HB2	2.02	0.42
1:D:95:LEU:HD23	1:D:95:LEU:HA	1.90	0.42
1:C:35:MET:HE2	1:C:60:PHE:HZ	1.85	0.42
1:B:137:VAL:HG13	1:B:146:ARG:HG2	2.01	0.42
1:D:28:CYS:O	1:D:32:VAL:HG23	2.20	0.41
1:D:9:GLN:HG2	1:D:11:THR:O	2.20	0.41
1:A:0:HIS:CG	1:A:1:MET:H	2.39	0.41
1:C:51:THR:HG22	1:C:138:THR:HG23	2.02	0.41
1:B:100:GLN:H	1:B:100:GLN:HG3	1.74	0.40
1:A:92:TYR:O	1:A:96:ARG:HG3	2.21	0.40
1:C:51:THR:HA	1:C:137:VAL:O	2.21	0.40
1:C:67:LEU:HD12	1:C:68:SER:H	1.87	0.40
1:D:3:HIS:NE2	1:D:21:ASP:HB2	2.37	0.40
1:D:40:LEU:HD23	1:D:40:LEU:HA	1.88	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	124/158 (78%)	121 (98%)	3 (2%)	0	100	100
1	B	113/158 (72%)	110 (97%)	3 (3%)	0	100	100
1	C	113/158 (72%)	111 (98%)	2 (2%)	0	100	100
1	D	113/158 (72%)	111 (98%)	2 (2%)	0	100	100
All	All	463/632 (73%)	453 (98%)	10 (2%)	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	113/136 (83%)	110 (97%)	3 (3%)	39	63
1	B	107/136 (79%)	107 (100%)	0	100	100
1	C	104/136 (76%)	104 (100%)	0	100	100
1	D	106/136 (78%)	105 (99%)	1 (1%)	70	84
All	All	430/544 (79%)	426 (99%)	4 (1%)	70	84

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	80	GLN
1	A	153	LEU
1	A	154	THR
1	D	143	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	44	ASN
1	B	44	ASN
1	B	56	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	80	GLN
1	D	44	ASN
1	D	100	GLN

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 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 ⓘ

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	128/158 (81%)	0.32	1 (0%) 82 78	28, 40, 66, 80	0
1	B	117/158 (74%)	0.18	0 100 100	30, 40, 54, 72	0
1	C	117/158 (74%)	0.61	11 (9%) 14 11	38, 56, 93, 106	0
1	D	117/158 (74%)	0.50	4 (3%) 48 38	38, 51, 71, 87	0
All	All	479/632 (75%)	0.40	16 (3%) 49 40	28, 48, 74, 106	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	149	SER	3.4
1	C	150	THR	3.0
1	C	1	MET	2.6
1	D	26	ASN	2.5
1	D	144	ASP	2.4
1	C	95	LEU	2.4
1	C	134	SER	2.2
1	C	63	ASP	2.2
1	C	142	SER	2.2
1	C	144	ASP	2.1
1	C	148	PHE	2.1
1	A	0	HIS	2.1
1	D	22	TYR	2.1
1	C	139	ILE	2.0
1	D	5	ILE	2.0
1	C	92	TYR	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 [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.