



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:23 PM UTC

PDB ID : 1CGE / pdb_00001cge
Title : CRYSTAL STRUCTURES OF RECOMBINANT 19-KDA HUMAN FI-BROBLAST COLLAGENASE COMPLEXED TO ITSELF
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Deposited on : 1994-02-03
Resolution : 1.90 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

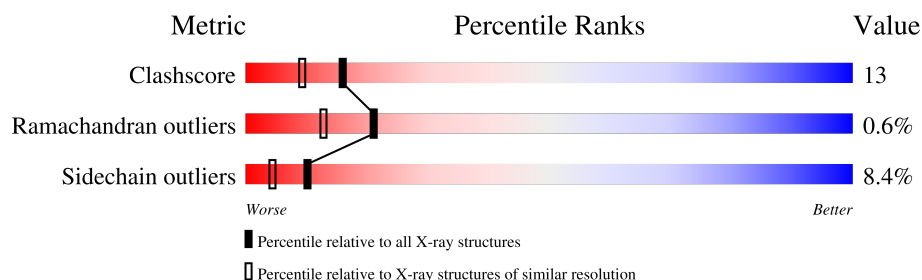
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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))
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	168	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 1853 atoms, of which 477 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 FIBROBLAST COLLAGENASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	162	Total	C	H	N	O	S	0	0	0
			1542	795	273	226	246	2			

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Zn	0	0
			2	2		

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

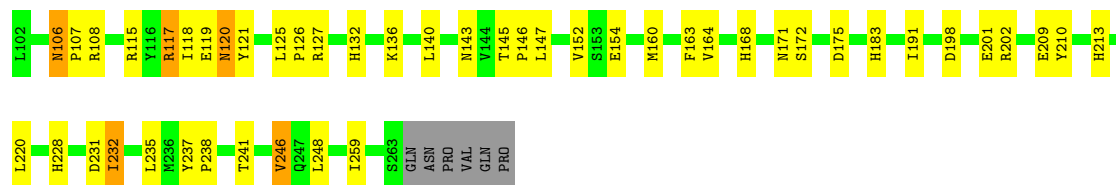
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	3	Total	Ca	0	0
			3	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	102	Total	H	O	0	0
			306	204	102		

Note EDS was not executed.

Chain A: 68% 25% . .



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	72.60Å 72.60Å 75.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	7.00 – 1.90	Depositor
% Data completeness (in resolution range)	(Not available) (7.00-1.90)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.198 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1853	wwPDB-VP
Average B, all atoms (Å ²)	16.0	wwPDB-VP

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, CA

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	0.59	0/1306	1.07	9/1779 (0.5%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	171	ASN	N-CA-C	6.57	120.21	112.72
1	A	237	TYR	N-CA-C	-6.07	102.12	109.83
1	A	209	GLU	N-CA-C	5.90	119.04	110.42
1	A	237	TYR	CA-C-N	5.84	125.38	119.19
1	A	237	TYR	C-N-CA	5.84	125.38	119.19
1	A	106	ASN	CA-C-N	5.62	125.95	119.93
1	A	106	ASN	C-N-CA	5.62	125.95	119.93
1	A	126	PRO	N-CA-C	-5.60	103.02	111.41
1	A	175	ASP	N-CA-C	5.19	119.16	112.41

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	1269	273	1158	32	0
2	A	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	3	0	0	0	0
4	A	102	204	0	6	0
All	All	1376	477	1158	32	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 13.

All (32) 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:A:202:ARG:HD3	4:A:601:HOH:O	1.83	0.79
1:A:107:PRO:O	1:A:108:ARG:HD3	1.97	0.65
1:A:120:ASN:HD21	1:A:163:PHE:H	1.47	0.62
1:A:235:LEU:HD22	1:A:248:LEU:HD23	1.81	0.62
1:A:136:LYS:HG2	4:A:590:HOH:O	1.99	0.61
1:A:106:ASN:HB2	1:A:108:ARG:CZ	2.32	0.59
1:A:106:ASN:HB2	1:A:108:ARG:NH2	2.18	0.58
1:A:120:ASN:N	1:A:120:ASN:HD22	2.03	0.57
1:A:213:HIS:CD2	1:A:246:VAL:HG11	2.39	0.56
1:A:121:TYR:CE2	1:A:127:ARG:HG3	2.43	0.54
1:A:164:VAL:HG21	1:A:168:HIS:CD2	2.44	0.53
1:A:228:HIS:CG	1:A:238:PRO:HG3	2.45	0.52
1:A:145:THR:HB	1:A:146:PRO:HD2	1.93	0.50
1:A:115:ARG:NH1	4:A:538:HOH:O	2.45	0.48
1:A:120:ASN:HD22	1:A:120:ASN:H	1.61	0.47
1:A:117:ARG:HG3	1:A:118:ILE:N	2.28	0.47
1:A:143:ASN:ND2	4:A:557:HOH:O	2.50	0.45
1:A:119:GLU:OE2	1:A:160:MET:HB3	2.17	0.45
1:A:108:ARG:HD2	1:A:259:ILE:HG21	1.98	0.44
1:A:117:ARG:NE	4:A:574:HOH:O	2.49	0.44
1:A:108:ARG:HD2	1:A:259:ILE:CG2	2.49	0.42
1:A:164:VAL:HG21	1:A:168:HIS:CG	2.54	0.42
1:A:235:LEU:HD11	1:A:246:VAL:HG23	2.02	0.42
1:A:232:ILE:O	1:A:232:ILE:HG13	2.20	0.42
1:A:117:ARG:NH2	4:A:574:HOH:O	2.51	0.42
1:A:140:LEU:HD11	1:A:246:VAL:HG22	2.02	0.41
1:A:172:SER:O	1:A:183:HIS:HE1	2.04	0.41
1:A:106:ASN:O	1:A:108:ARG:HG2	2.20	0.41
1:A:107:PRO:C	1:A:108:ARG:HD3	2.45	0.41
1:A:132:HIS:O	1:A:136:LYS:HB2	2.21	0.41
1:A:198:ASP:HB3	1:A:201:GLU:HG2	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:115:ARG:CZ	1:A:152:VAL:HG11	2.51	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	160/168 (95%)	157 (98%)	2 (1%)	1 (1%)	21 13

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	210	TYR

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	131/140 (94%)	120 (92%)	11 (8%)	10 4

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	117	ARG
1	A	120	ASN

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Mol	Chain	Res	Type
1	A	125	LEU
1	A	147	LEU
1	A	154	GLU
1	A	191	ILE
1	A	220	LEU
1	A	231	ASP
1	A	232	ILE
1	A	241	THR
1	A	246	VAL

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	106	ASN
1	A	120	ASN
1	A	186	GLN
1	A	213	HIS
1	A	247	GLN
1	A	257	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](#)

Of 5 ligands modelled in this entry, 5 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.