

Full wwPDB X-ray Structure Validation Report (i)

Aug 20, 2023 – 02:02 PM EDT

PDB ID	:	2HZ9
Title	:	Crystal structure of Lys12Val/Asn95Val/Cys117Val mutant of human acidic
		fibroblast growth factor at 1.70 angstrom resolution.
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Deposited on		
Resolution	:	1.70 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 (i) 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 (1)) were used in the production of this report:

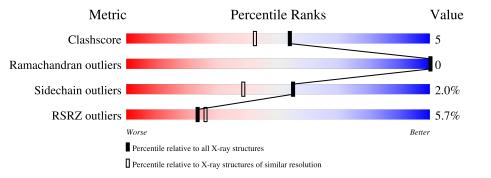
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	$\begin{array}{c} \textbf{Whole archive} \\ (\# \textbf{Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$		
Clashscore	141614	4695 (1.70-1.70)		
Ramachandran outliers	138981	4610 (1.70-1.70)		
Sidechain outliers	138945	4610 (1.70-1.70)		
RSRZ outliers	127900	4222 (1.70-1.70)		

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 <=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	А	146	88%	6% • •					
1	В	146	<mark>6%</mark> 87%	10% • •					



2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 2526 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.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	۸	141	Total	С	Ν	0	S	0	0	0
	I A	141	1134	722	201	208	3	0		
1	р	142	Total	С	Ν	0	S	0	0	0
	D	142	1144	728	204	209	3	0		0

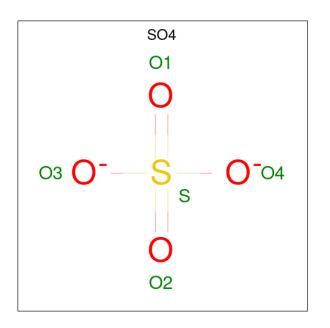
• Molecule 1 is a protein called Heparin-binding growth factor 1.

Chain	Residue	Modelled	Actual	Comment	Reference
A	1C	HIS	-	- expression tag	
А	1C	HIS	-	expression tag	UNP P05230
A	1C	HIS	-	expression tag	UNP P05230
А	1D	HIS	-	expression tag	UNP P05230
A	1E	HIS	-	expression tag	UNP P05230
А	1F	HIS	-	expression tag	UNP P05230
A	12	VAL	LYS	engineered mutation	UNP P05230
А	95	VAL	ASN	engineered mutation	UNP P05230
А	117	VAL	CYS	engineered mutation	UNP P05230
В	1B	HIS	-	expression tag	UNP P05230
В	1B	HIS	-	expression tag	UNP P05230
В	1C	HIS	-	expression tag	UNP P05230
В	1D	HIS	-	expression tag	UNP P05230
В	1E	HIS	-	expression tag	UNP P05230
В	1F	HIS	-	expression tag	UNP P05230
В	12	VAL	LYS	engineered mutation	UNP P05230
В	95	VAL	ASN engineered mutation		UNP P05230
В	117	VAL	CYS	engineered mutation	UNP P05230

There are 18 discrepancies between the modelled and reference sequences:

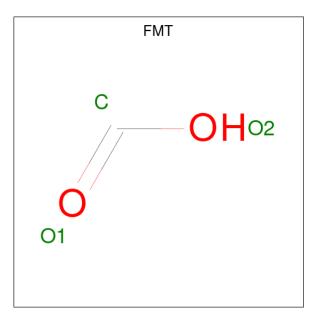
• Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O_4S).





Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	А	1	Total 5	0 4	S 1	0	0

• Molecule 3 is FORMIC ACID (three-letter code: FMT) (formula: CH_2O_2).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 3 & 1 & 2 \end{array}$	0	0
3	В	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 3 & 1 & 2 \end{array}$	0	0
3	В	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 3 & 1 & 2 \end{array}$	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	В	1	Total 3	C 1	0 2	0	0

• Molecule 4 is water.

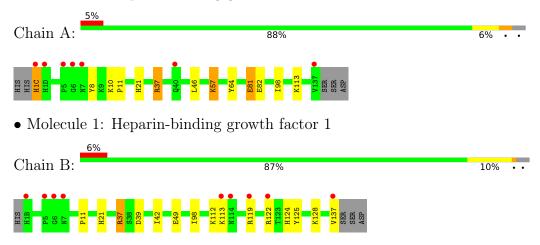
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	119	Total O 119 119	0	0
4	В	112	Total O 112 112	0	0



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: Heparin-binding growth factor 1





4 Data and refinement statistics (i)

Property	Value	Source	
Space group	C 2 2 21	Depositor	
Cell constants	76.28Å 95.37Å 107.33Å	Deperitor	
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor	
Resolution (Å)	43.58 - 1.70	Depositor	
Resolution (A)	43.58 - 1.70	EDS	
% Data completeness	(Not available) $(43.58-1.70)$	Depositor	
(in resolution range)	96.4(43.58-1.70)	EDS	
R _{merge}	0.08	Depositor	
R _{sym}	(Not available)	Depositor	
$< I/\sigma(I) > 1$	$1.12 (at 1.70 \text{\AA})$	Xtriage	
Refinement program	CNS	Depositor	
D D.	0.199 , 0.226	Depositor	
R, R_{free}	0.194 , (Not available)	DCC	
R_{free} test set	No test flags present.	wwPDB-VP	
Wilson B-factor $(Å^2)$	23.7	Xtriage	
Anisotropy	0.483	Xtriage	
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.36 , 42.0	EDS	
L-test for twinning ²	$ \langle L \rangle = 0.49, \langle L^2 \rangle = 0.33$	Xtriage	
Estimated twinning fraction	No twinning to report.	Xtriage	
F_o, F_c correlation	0.96	EDS	
Total number of atoms	2526	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 7.32% of the height of the origin peak. No significant pseudotranslation is detected.

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.



¹Intensities estimated from amplitudes.

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: FMT, SO4

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		
	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.65	0/1165	0.78	0/1577	
1	В	0.59	0/1176	0.74	1/1592~(0.1%)	
All	All	0.62	0/2341	0.76	1/3169~(0.0%)	

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	В	49	GLU	N-CA-C	-5.15	97.11	111.00

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	А	1134	0	1112	12	1
1	В	1144	0	1119	12	0
2	А	5	0	0	0	0
3	А	3	0	1	0	0
3	В	9	0	3	1	0
4	А	119	0	0	5	0
4	В	112	0	0	2	0
All	All	2526	0	2235	24	1



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 5.

All (24) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:B:11:PRO:HG2	1:B:137:VAL:HG21	1.65	0.79
1:B:11:PRO:HG2	1:B:137:VAL:CG2	2.25	0.66
1:A:81:GLU:HG2	1:A:82:GLU:N	2.13	0.64
1:B:112:LYS:NZ	1:B:122:ARG:HH22	1.97	0.63
1:A:10:LYS:HB2	1:A:11:PRO:HD2	1.81	0.62
1:A:57:LYS:HD2	1:A:64:TYR:CE2	2.37	0.59
1:A:57:LYS:NZ	4:A:229:HOH:O	2.36	0.58
1:B:37:ARG:HD3	4:B:206:HOH:O	2.07	0.54
1:A:98:ILE:HG13	4:A:168:HOH:O	2.10	0.52
1:B:112:LYS:HZ3	1:B:122:ARG:HH22	1.57	0.50
1:A:37:ARG:NH1	1:A:37:ARG:HG2	2.28	0.49
1:A:98:ILE:CD1	4:A:168:HOH:O	2.61	0.48
1:A:37:ARG:HG2	1:A:37:ARG:HH11	1.79	0.47
1:B:112:LYS:CE	1:B:122:ARG:HH22	2.27	0.47
1:B:98:ILE:HG13	4:B:153:HOH:O	2.16	0.46
1:B:21:HIS:HE1	1:B:113:LYS:O	2.00	0.45
1:A:1(C):HIS:HB2	4:A:256:HOH:O	2.16	0.44
1:A:8:TYR:HB3	1:A:46:LEU:O	2.18	0.43
1:B:128:LYS:H	3:B:143:FMT:C	2.33	0.41
1:A:21:HIS:HE1	1:A:113:LYS:O	2.04	0.41
1:B:124:HIS:HD2	1:B:125:TYR:O	2.03	0.41
1:B:119:ARG:NH2	1:B:122:ARG:HD2	2.36	0.41
1:B:39:ASP:O	1:B:42:ILE:HG12	2.22	0.40
1:A:57:LYS:CG	4:A:192:HOH:O	2.70	0.40

All (1) 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:A:81:GLU:OE2	1:A:81:GLU:OE2[4_566]	2.03	0.17



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	Perce	ntiles
1	А	139/146~(95%)	135~(97%)	4(3%)	0	100	100
1	В	140/146~(96%)	132 (94%)	8 (6%)	0	100	100
All	All	279/292~(96%)	267~(96%)	12~(4%)	0	100	100

There are no Ramachandran outliers to report.

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	Percent	iles
1	А	124/129~(96%)	120~(97%)	4(3%)	39 2	0
1	В	125/129~(97%)	124 (99%)	1 (1%)	81 7	4
All	All	249/258~(96%)	244~(98%)	5(2%)	55 3	8

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	1(C)	HIS
1	А	37	ARG
1	А	57	LYS
1	А	81	GLU
1	В	37	ARG

Sometimes side chains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such side chains are listed below:



Mol	Chain	Res	Type
1	А	77	GLN
1	В	1(B)	HIS
1	В	21	HIS
1	В	124	HIS

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)

5 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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| > 2 is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Trune	Chain			B	Bond lengths			Bond angles		
10101	Type	Unam	Res	Link	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z >2	
3	FMT	В	143	-	2,2,2	0.73	0	$1,\!1,\!1$	0.61	0	
3	FMT	А	142	-	2,2,2	1.10	0	$1,\!1,\!1$	0.56	0	
3	FMT	В	142	-	2,2,2	0.89	0	$1,\!1,\!1$	0.63	0	
3	FMT	В	141	-	2,2,2	0.83	0	$1,\!1,\!1$	0.62	0	
2	SO4	А	141	-	4,4,4	0.29	0	6,6,6	0.33	0	

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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	В	143	FMT	1	0

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, 95^{th} 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	$\langle RSRZ \rangle$	# RSRZ > 2	$OWAB(Å^2)$	Q<0.9
1	А	141/146~(96%)	0.19	7 (4%) 28 32	16, 24, 40, 54	0
1	В	142/146~(97%)	0.33	9 (6%) 20 22	17, 27, 45, 55	0
All	All	283/292~(96%)	0.26	16 (5%) 23 26	16, 25, 44, 55	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	6	GLY	5.7
1	А	6	GLY	4.8
1	А	7	ASN	4.7
1	В	137	VAL	3.7
1	В	1(B)	HIS	3.6
1	А	1(C)	HIS	3.3
1	А	5	PRO	3.3
1	В	7	ASN	2.7
1	А	1(D)	HIS	2.7
1	В	119	ARG	2.6
1	В	113	LYS	2.6
1	А	137	VAL	2.5
1	В	122	ARG	2.3
1	В	5	PRO	2.3
1	А	40	GLN	2.3
1	В	114	ASN	2.2

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)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	FMT	А	142	3/3	0.76	0.22	37,37,42,44	0
3	FMT	В	142	3/3	0.76	0.15	45,45,47,47	0
3	FMT	В	141	3/3	0.86	0.11	55,55,57,57	0
3	FMT	В	143	3/3	0.86	0.16	36,36,40,41	0
2	SO4	А	141	5/5	0.96	0.13	37,38,41,41	0

6.5 Other polymers (i)

There are no such residues in this entry.

