



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 17, 2021 – 02:52 AM EDT

PDB ID : 1JT4
Title : Human Acidic Fibroblast Growth Factor. 141 Amino Acid Form with Amino Terminal His Tag AND VAL 109 REPLACED BY LEU (V109L)
Authors : Brych, S.R.; Blaber, S.I.; Logan, T.M.; Blaber, M.
Deposited on : 2001-08-20
Resolution : 1.78 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 following versions of software and data (see [references ⓘ](#)) 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 : 2.23.2
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) : 2.23.2

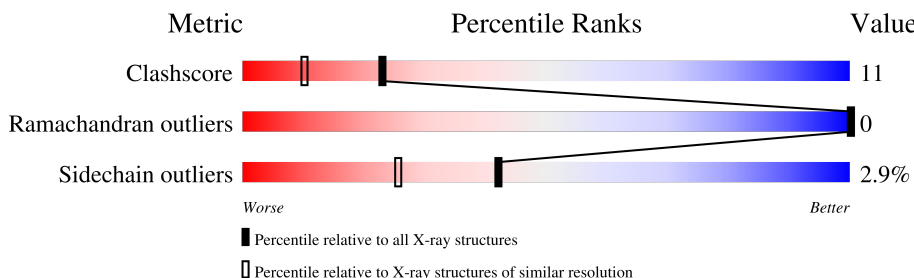
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.78 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	141614	10184 (1.80-1.76)
Ramachandran outliers	138981	10051 (1.80-1.76)
Sidechain outliers	138945	10050 (1.80-1.76)

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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	146	
1	B	146	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2480 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 acidic fibroblast growth factor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	141	Total	C	N	O	S	0	0	0
			1131	718	200	209	4			
1	B	141	Total	C	N	O	S	0	0	0
			1129	717	202	206	4			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1A	HIS	-	expression tag	UNP P05230
A	1B	HIS	-	expression tag	UNP P05230
A	1C	HIS	-	expression tag	UNP P05230
A	1D	HIS	-	expression tag	UNP P05230
A	1E	HIS	-	expression tag	UNP P05230
A	1F	HIS	-	expression tag	UNP P05230
A	109	LEU	VAL	engineered mutation	UNP P05230
B	1A	HIS	-	expression tag	UNP P05230
B	1B	HIS	-	expression tag	UNP P05230
B	1C	HIS	-	expression tag	UNP P05230
B	1D	HIS	-	expression tag	UNP P05230
B	1E	HIS	-	expression tag	UNP P05230
B	1F	HIS	-	expression tag	UNP P05230
B	109	LEU	VAL	engineered mutation	UNP P05230

- Molecule 2 is FORMIC ACID (three-letter code: FMT) (formula: CH₂O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			3	1	2		
2	A	1	Total	C	O	0	0
			3	1	2		
2	B	1	Total	C	O	0	0
			3	1	2		
2	B	1	Total	C	O	0	0
			3	1	2		

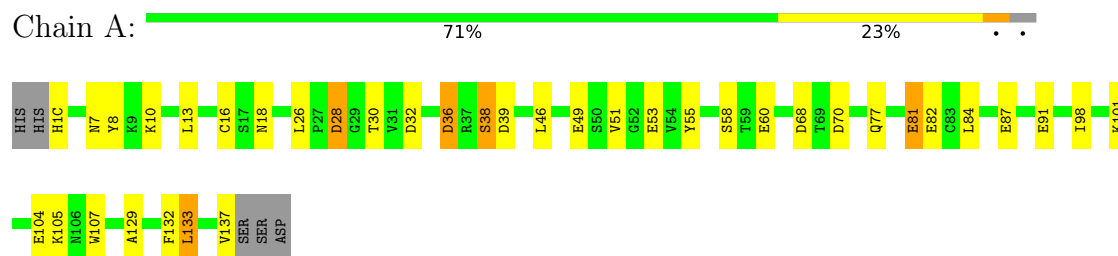
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	108	Total	O	0	0
			108	108		
3	B	100	Total	O	0	0
			100	100		

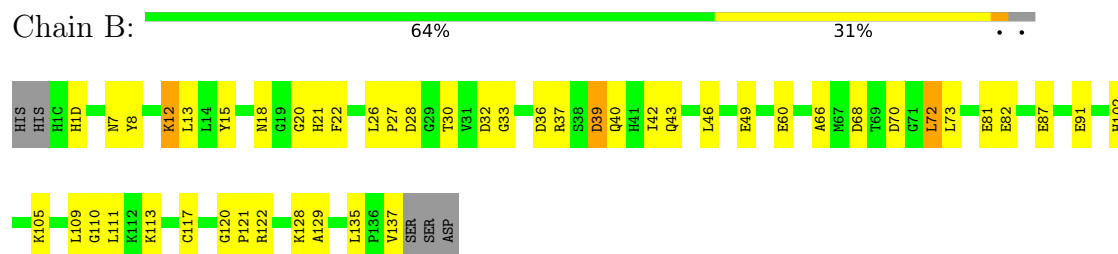
3 Residue-property plots [i](#)

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. 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. 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: acidic fibroblast growth factor



- Molecule 1: acidic fibroblast growth factor



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	73.56Å 97.36Å 109.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.90 – 1.78 44.45 – 1.78	Depositor EDS
% Data completeness (in resolution range)	84.3 (39.90-1.78) 93.1 (44.45-1.78)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.80 (at 1.77Å)	Xtriage
Refinement program	TNT	Depositor
R, R_{free}	(Not available) , (Not available) 0.187 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	19.2	Xtriage
Anisotropy	0.340	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 75.5	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	2480	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.43% 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

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

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	0.85	8/1162 (0.7%)	1.05	10/1571 (0.6%)
1	B	0.83	6/1160 (0.5%)	1.07	12/1568 (0.8%)
All	All	0.84	14/2322 (0.6%)	1.06	22/3139 (0.7%)

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	104	GLU	CD-OE2	6.23	1.32	1.25
1	B	87	GLU	CD-OE2	5.95	1.32	1.25
1	A	60	GLU	CD-OE2	5.72	1.31	1.25
1	A	91	GLU	CD-OE2	5.68	1.31	1.25
1	A	49	GLU	CD-OE2	5.64	1.31	1.25
1	B	81	GLU	CD-OE2	5.64	1.31	1.25
1	B	82	GLU	CD-OE2	5.60	1.31	1.25
1	B	91	GLU	CD-OE2	5.56	1.31	1.25
1	B	60	GLU	CD-OE2	5.49	1.31	1.25
1	A	53	GLU	CD-OE2	5.49	1.31	1.25
1	B	49	GLU	CD-OE1	5.47	1.31	1.25
1	A	81	GLU	CD-OE2	5.15	1.31	1.25
1	A	82	GLU	CD-OE2	5.11	1.31	1.25
1	A	87	GLU	CD-OE2	5.09	1.31	1.25

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	68	ASP	CB-CG-OD1	7.22	124.80	118.30
1	A	68	ASP	CB-CG-OD2	-7.20	111.82	118.30
1	A	39	ASP	CB-CG-OD2	-7.00	112.00	118.30
1	B	36	ASP	CB-CG-OD1	6.96	124.57	118.30
1	A	36	ASP	CB-CG-OD1	6.94	124.55	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	36	ASP	CB-CG-OD2	-6.88	112.11	118.30
1	A	36	ASP	CB-CG-OD2	-6.75	112.22	118.30
1	A	32	ASP	CB-CG-OD1	6.48	124.13	118.30
1	B	68	ASP	CB-CG-OD2	-6.48	112.47	118.30
1	B	70	ASP	CB-CG-OD2	-6.33	112.60	118.30
1	B	28	ASP	CB-CG-OD2	-6.24	112.69	118.30
1	A	32	ASP	CB-CG-OD2	-6.16	112.76	118.30
1	B	32	ASP	CB-CG-OD2	-6.07	112.83	118.30
1	B	39	ASP	CB-CG-OD2	-5.99	112.91	118.30
1	B	68	ASP	CB-CG-OD1	5.89	123.60	118.30
1	A	28	ASP	CB-CG-OD2	-5.61	113.26	118.30
1	B	28	ASP	CB-CG-OD1	5.48	123.23	118.30
1	B	70	ASP	CB-CG-OD1	5.41	123.16	118.30
1	B	32	ASP	CB-CG-OD1	5.33	123.10	118.30
1	A	39	ASP	CB-CG-OD1	5.29	123.06	118.30
1	B	39	ASP	CB-CG-OD1	5.09	122.89	118.30
1	A	70	ASP	CB-CG-OD2	-5.04	113.76	118.30

There are no chirality outliers.

There are no planarity outliers.

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	1131	0	1100	21	0
1	B	1129	0	1101	30	0
2	A	6	0	2	1	0
2	B	6	0	2	2	0
3	A	108	0	0	5	0
3	B	100	0	0	3	1
All	All	2480	0	2205	51	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 11.

All (51) 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:26:LEU:HB2	1:B:30:THR:HG22	1.29	1.13
1:A:13:LEU:CD2	1:A:137:VAL:HG12	2.11	0.80
1:B:26:LEU:HD12	1:B:30:THR:CG2	2.16	0.74
1:B:26:LEU:HB2	1:B:30:THR:CG2	2.18	0.66
1:B:137:VAL:HG22	3:B:1261:HOH:O	1.96	0.66
1:B:18:ASN:HB2	1:B:129:ALA:HA	1.79	0.65
1:B:26:LEU:HB3	1:B:27:PRO:HD2	1.79	0.65
1:B:72:LEU:N	1:B:72:LEU:HD23	2.12	0.64
1:B:7:ASN:HB2	3:B:1250:HOH:O	1.96	0.63
1:B:26:LEU:HD12	1:B:30:THR:HG21	1.80	0.63
1:A:13:LEU:HD21	1:A:137:VAL:HG12	1.81	0.62
1:A:28:ASP:OD1	1:A:30:THR:HB	2.02	0.59
1:A:10:LYS:HB2	1:A:10:LYS:NZ	2.21	0.56
1:B:128:LYS:H	2:B:163:FMT:C	2.18	0.55
1:A:81:GLU:HG2	3:A:337:HOH:O	2.06	0.55
1:B:12:LYS:HA	1:B:137:VAL:HG13	1.87	0.54
1:B:102:HIS:HB3	1:B:105:LYS:HD2	1.89	0.54
1:A:105:LYS:HE3	1:A:107:TRP:CZ2	2.44	0.53
1:B:120:GLY:N	1:B:121:PRO:HD2	2.25	0.52
1:B:109:LEU:HD12	1:B:110:GLY:N	2.24	0.52
1:A:98:ILE:HD11	3:A:278:HOH:O	2.09	0.51
1:A:77:GLN:NE2	3:A:334:HOH:O	2.43	0.51
1:B:12:LYS:HB2	1:B:135:LEU:O	2.10	0.51
1:B:39:ASP:O	1:B:43:GLN:NE2	2.40	0.50
1:A:13:LEU:HD23	1:A:137:VAL:HG12	1.92	0.50
1:B:13:LEU:CD2	1:B:137:VAL:HG12	2.41	0.49
1:B:21:HIS:HE1	1:B:113:LYS:O	1.96	0.48
1:B:40:GLN:O	1:B:43:GLN:OE1	2.33	0.47
1:A:1(C):HIS:HB3	2:A:160:FMT:O1	2.14	0.47
1:B:111:LEU:HD23	1:B:117:CYS:HA	1.97	0.47
1:A:8:TYR:HB3	1:A:46:LEU:O	2.16	0.46
1:B:40:GLN:HA	1:B:43:GLN:HE22	1.81	0.46
1:A:16:CYS:HB2	1:A:132:PHE:CE2	2.52	0.45
1:B:66:ALA:O	1:B:73:LEU:HA	2.17	0.45
1:B:120:GLY:N	1:B:121:PRO:CD	2.80	0.45
1:B:15:TYR:CZ	1:B:20:GLY:HA2	2.52	0.45
1:B:72:LEU:N	1:B:72:LEU:CD2	2.79	0.45
1:B:137:VAL:N	3:B:1261:HOH:O	2.28	0.44
1:A:10:LYS:NZ	1:A:10:LYS:CB	2.80	0.44
1:A:7:ASN:HB2	3:A:250:HOH:O	2.19	0.43
1:A:81:GLU:OE2	1:A:101:LYS:HE2	2.19	0.43
1:B:1(D):HIS:H	2:B:162:FMT:C	2.31	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:26:LEU:HD12	1:A:30:THR:HG22	2.00	0.42
1:B:22:PHE:O	1:B:33:GLY:HA2	2.19	0.42
1:A:133:LEU:HD12	1:A:133:LEU:HA	1.89	0.42
1:B:8:TYR:HB3	1:B:46:LEU:O	2.21	0.41
1:A:55:TYR:CZ	1:A:84:LEU:HD21	2.55	0.41
1:A:36:ASP:OD1	1:A:38:SER:N	2.44	0.41
1:A:51:VAL:HG23	3:A:260:HOH:O	2.20	0.41
1:B:39:ASP:O	1:B:42:ILE:HG12	2.21	0.41
1:A:18:ASN:HB2	1:A:129:ALA:HA	2.03	0.41

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)
3:B:1403:HOH:O	3:B:1419:HOH:O[3_556]	2.06	0.14

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	139/146 (95%)	133 (96%)	6 (4%)	0	100	100
1	B	139/146 (95%)	134 (96%)	5 (4%)	0	100	100
All	All	278/292 (95%)	267 (96%)	11 (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	Percentiles	
1	A	123/129 (95%)	120 (98%)	3 (2%)	49	33
1	B	122/129 (95%)	118 (97%)	4 (3%)	38	21
All	All	245/258 (95%)	238 (97%)	7 (3%)	42	25

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	SER
1	A	58	SER
1	A	133	LEU
1	B	12	LYS
1	B	37	ARG
1	B	72	LEU
1	B	122	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	77	GLN
1	B	77	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 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	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	FMT	B	162	-	0,2,2	-	-	0,1,1	-	-
2	FMT	B	163	-	0,2,2	-	-	0,1,1	-	-
2	FMT	A	161	-	0,2,2	-	-	0,1,1	-	-
2	FMT	A	160	-	0,2,2	-	-	0,1,1	-	-

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.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	162	FMT	1	0
2	B	163	FMT	1	0
2	A	160	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](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.