



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 1HUW / pdb_00001huw
Title : THE CRYSTAL STRUCTURE OF AFFINITY-MATURED HUMAN GROWTH HORMONE AT 2 ANGSTROMS RESOLUTION
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Deposited on : 1993-09-22
Resolution : 2.00 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

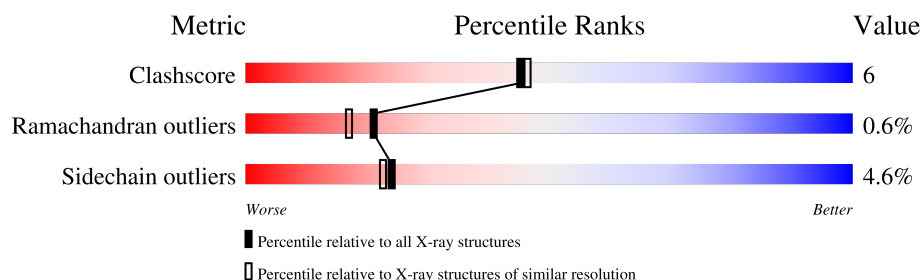
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)

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	191	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1438 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 HUMAN GROWTH HORMONE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	166	Total	C	N	O	S	0	3	0
			1361	870	223	262	6			

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	10	ALA	PHE	engineered mutation	UNP P01241
A	14	TRP	MET	engineered mutation	UNP P01241
A	18	ASP	HIS	engineered mutation	UNP P01241
A	21	ASN	HIS	engineered mutation	UNP P01241
A	41	ILE	LYS	engineered mutation	UNP P01241
A	42	HIS	TYR	engineered mutation	UNP P01241
A	45	TRP	LEU	engineered mutation	UNP P01241
A	46	TRP	GLN	engineered mutation	UNP P01241
A	54	PRO	PHE	engineered mutation	UNP P01241
A	64	LYS	ARG	engineered mutation	UNP P01241
A	167	ASN	ARG	engineered mutation	UNP P01241
A	171	SER	ASP	engineered mutation	UNP P01241
A	174	SER	GLU	engineered mutation	UNP P01241
A	176	TYR	PHE	engineered mutation	UNP P01241
A	179	THR	ILE	engineered mutation	UNP P01241

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	77	Total	O	0	0
			77	77		

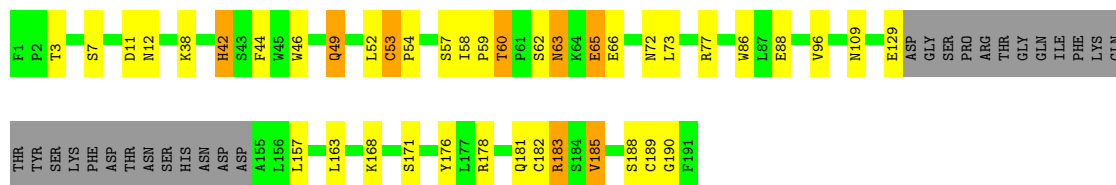
3 Residue-property plots [i](#)

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.

Note EDS was not executed.

• Molecule 1: HUMAN GROWTH HORMONE

Chain A: 



4 Data and refinement statistics

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

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	80.45Å 59.34Å 50.87Å 90.00° 128.20° 90.00°	Depositor
Resolution (Å)	8.00 – 2.00	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-2.00)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.185 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1438	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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	1.03	2/1403 (0.1%)	1.74	18/1903 (0.9%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	42	HIS	CD2-NE2	-6.16	1.31	1.37
1	A	60	THR	CA-CB	5.23	1.60	1.54

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	57	SER	N-CA-C	7.40	120.40	111.82
1	A	3	THR	N-CA-CB	-6.59	100.94	110.56
1	A	53	CYS	CA-C-N	6.16	125.84	119.56
1	A	53	CYS	C-N-CA	6.16	125.84	119.56
1	A	62	SER	N-CA-C	5.84	117.72	111.36
1	A	185	VAL	CA-C-N	-5.83	114.08	123.23
1	A	185	VAL	C-N-CA	-5.83	114.08	123.23
1	A	183	ARG	CB-CG-CD	-5.82	97.91	111.30
1	A	86	TRP	CE2-CD2-CG	-5.67	100.39	107.20
1	A	66	GLU	CA-CB-CG	-5.65	102.81	114.10
1	A	42	HIS	CB-CG-CD2	-5.34	124.26	131.20
1	A	86	TRP	CD1-CG-CD2	5.25	114.70	106.30
1	A	88	GLU	CA-C-N	5.18	126.31	119.84
1	A	88	GLU	C-N-CA	5.18	126.31	119.84
1	A	86	TRP	CG-CD1-NE1	-5.13	103.53	110.20
1	A	49	GLN	OE1-CD-NE2	-5.11	117.49	122.60
1	A	109	ASN	N-CA-C	-5.07	98.32	107.75
1	A	12	ASN	O-C-N	5.07	127.92	122.15

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	1361	0	1328	16	0
2	A	77	0	0	0	0
All	All	1438	0	1328	16	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 6.

All (16) 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:7:SER:HB3	1:A:185:VAL:HB	1.45	0.98
1:A:44:PHE:HE2	1:A:49:GLN:HG2	1.46	0.78
1:A:38:LYS:HE2	1:A:42:HIS:HE1	1.61	0.65
1:A:49:GLN:HG3	1:A:52:LEU:HG	1.86	0.58
1:A:72:ASN:OD1	1:A:183:ARG:HD2	2.07	0.54
1:A:73:LEU:HD21	1:A:129:GLU:HG2	1.89	0.54
1:A:63:ASN:OD1	1:A:65:GLU:HG2	2.08	0.53
1:A:182:CYS:SG	1:A:190:GLY:N	2.88	0.46
1:A:96:VAL:HG11	1:A:163:LEU:HD11	1.98	0.46
1:A:53:CYS:SG	1:A:168:LYS:NZ	2.90	0.45
1:A:7:SER:HB3	1:A:185:VAL:CB	2.32	0.45
1:A:53:CYS:HA	1:A:54:PRO:HD3	1.85	0.42
1:A:42:HIS:O	1:A:46:TRP:HD1	2.03	0.42
1:A:178:ARG:HD2	1:A:181:GLN:OE1	2.20	0.42
1:A:60:THR:HB	1:A:176:TYR:OH	2.21	0.41
1:A:58:ILE:HA	1:A:59:PRO:HD2	1.97	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	165/191 (86%)	160 (97%)	4 (2%)	1 (1%)	21 17

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	188	SER

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	155/175 (89%)	148 (96%)	7 (4%)	24 23

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

Mol	Chain	Res	Type
1	A	11	ASP
1	A	63	ASN
1	A	65	GLU
1	A	77	ARG
1	A	157	LEU
1	A	171	SER
1	A	189	CYS

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

Mol	Chain	Res	Type
1	A	40	GLN
1	A	42	HIS
1	A	122	GLN
1	A	181	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

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.