

wwPDB X-ray Structure Validation Summary Report (i)

May 30, 2020 – 04:49 am BST

PDB ID : 1QIY

Title : HUMAN INSULIN HEXAMERS WITH CHAIN B HIS MUTATED TO TYR

COMPLEXED WITH PHENOL

Authors: Tang, L.; Whittingham, J.L.; Verma, C.S.; Caves, L.S.D.; Dodson, G.G.

Deposited on : 1999-06-18

Resolution : 2.30 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 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) : NOT EXECUTED EDS : NOT EXECUTED

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

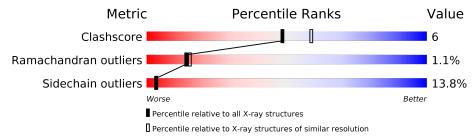
Validation Pipeline (wwPDB-VP) : 2.11

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 2.30 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	Similar resolution
Metric	$(\# { m Entries})$	$(\# ext{Entries}, ext{resolution range}(ext{Å}))$
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)

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 on 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%

Note EDS was not executed.





2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 2596 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 INSULIN A CHAIN.

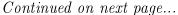
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
1	A	21	Total C N O S 163 99 25 35 4	12	0	0
1	C	21	Total C N O S	2	0	0
1	E	21	163 99 25 35 4 Total C N O S	5	0	0
1			163 99 25 35 4 Total C N O S			
1	G	21	163 99 25 35 4	5	0	0
1	I	21	Total C N O S 170 105 25 36 4	0	1	0
1	K	21	Total C N O S 163 99 25 35 4	0	0	0

• Molecule 2 is a protein called INSULIN B CHAIN.

Mol	Chain	Residues		Ato	ms			ZeroOcc	AltConf	Trace	
2	В	30	Total	С	N	О	S	14	0	0	
	Ъ	30	244	161	38	43	2	14	0	0	
2	D	30	Total	С	N	Ο	S	24	2	0	
	D	30	254	169	38	45	2	24	<u> </u>	U	
2	F	30	Total	С	N	Ο	S	35	0	0	
	I.	30	244	161	38	43	2	30			
2	Н	30	Total	С	N	О	S	27	0	0	
	11	30	244	161	38	43	2	21	U		
2	J	30	Total	С	N	О	S	21	0	0	
	9	30	244	161	38	43	2	21	0	0	
2	L	30	Total	С	N	О	S	17	0	0	
		30	244	161	38	43	2	11	U	0	

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
В	5	TYR	HIS	engineered mutation	UNP P01308

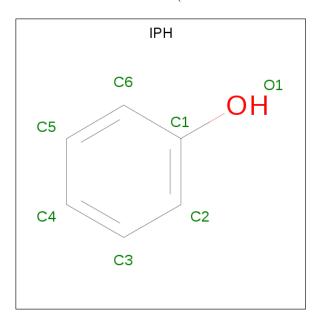




 $Continued\ from\ previous\ page...$

Chain	Residue	Modelled	Actual	Comment	Reference
D	5	TYR	HIS	engineered mutation	UNP P01308
F	5	TYR	HIS	engineered mutation	UNP P01308
Н	5	TYR	HIS	engineered mutation	UNP P01308
J	5	TYR	HIS	engineered mutation	UNP P01308
L	5	TYR	HIS	engineered mutation	UNP P01308

• Molecule 3 is PHENOL (three-letter code: IPH) (formula: C₆H₆O).



Mol	Chain	Residues	${f Atoms}$	ZeroOcc	AltConf
3	A	1	Total C O 7 6 1	0	0
3	С	1	Total C O 7 6 1	0	0
3	Е	1	Total C O 7 6 1	0	0
3	G	1	Total C O 7 6 1	0	0
3	I	1	Total C O 7 6 1	0	0
3	K	1	Total C O 7 6 1	0	0

• Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	J	1	Total Zn 1 1	0	0
4	L	1	Total Zn 1 1	0	0

 \bullet Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	J	1	Total Cl 1 1	0	0
5	L	1	Total Cl 1 1	0	0

• Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	4	Total O 4 4	0	0
6	В	9	Total O 9 9	0	0
6	С	9	Total O 9 9	0	0
6	D	10	Total O 10 10	0	0
6	Е	3	Total O 3 3	0	0
6	F	6	Total O 6 6	0	0
6	G	4	Total O 4 4	0	0
6	Н	10	Total O 10 10	0	0
6	I	7	Total O 7 7	0	0
6	J	13	Total O 13 13	0	0
6	K	8	Total O 8 8	0	0
6	L	8	Total O 8 8	0	0

 ${\tt SEQUENCE-PLOTS\ INFOmissing INFO}$



3 Data and refinement statistics (i)

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants	61.10Å 62.08Å 48.35Å	Depositor
a, b, c, α , β , γ	90.00° 109.87° 90.00°	Depositor
Resolution (Å)	36.80 - 2.30	Depositor
% Data completeness	92.0 (36.80-2.30)	Depositor
(in resolution range)	32.0 (30.00 2.30)	Depositor
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROLSQ	Depositor
R, R_{free}	0.186 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2596	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP



4 Model quality (i)

4.1 Standard geometry (i)

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

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 Bo		nd lengths	Во	ond angles
MIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5
1	A	0.58	0/164	1.39	1/220~(0.5%)
1	С	0.72	0/164	1.55	$3/220 \ (1.4\%)$
1	Е	0.56	0/164	1.24	0/220
1	G	0.85	0/164	1.50	1/220~(0.5%)
1	I	0.63	0/177	1.42	$1/238 \ (0.4\%)$
1	K	0.59	0/164	1.42	$2/220 \ (0.9\%)$
2	В	0.80	$1/251 \ (0.4\%)$	1.42	5/338 (1.5%)
2	D	1.07	2/272~(0.7%)	1.78	11/366 (3.0%)
2	F	0.95	$1/251 \ (0.4\%)$	1.77	6/338 (1.8%)
2	Н	1.14	$2/251 \ (0.8\%)$	2.48	5/338 (1.5%)
2	J	1.91	$2/251 \; (0.8\%)$	1.92	10/338 (3.0%)
2	L	0.71	$1/251 \; (0.4\%)$	1.91	10/338 (3.0%)
All	All	0.99	9/2524~(0.4%)	1.73	55/3394 (1.6%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	D	1	0
2	F	1	0
2	Н	0	1
2	L	0	1
All	All	2	2

The worst 5 of 9 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\operatorname{\AA})$	$\operatorname{Ideal}(ext{\AA})$
2	J	30	THR	C-OXT	-27.30	0.71	1.23
2	D	30	THR	C-OXT	11.59	1.45	1.23

Continued on next page...



 $Continued\ from\ previous\ page...$

Mol	Chain	Res	Type	Atoms	${f Z}$	${ m Observed}({ m \AA})$	$\operatorname{Ideal}(ext{\AA})$
2	Н	29	LYS	C-N	11.36	1.60	1.34
2	F	29	LYS	CA-CB	-9.43	1.33	1.53
2	Н	29	LYS	CA-CB	-8.55	1.35	1.53

The worst 5 of 55 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(^o)$	$\mathbf{Ideal}(^o)$
2	Н	22	ARG	CD-NE-CZ	24.09	157.33	123.60
2	Н	29	LYS	O-C-N	-21.36	88.53	122.70
2	Н	22	ARG	NE-CZ-NH2	12.62	126.61	120.30
2	Н	29	LYS	N-CA-CB	11.76	131.77	110.60
2	L	22	ARG	NE-CZ-NH2	-11.52	114.54	120.30

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	D	1	PHE	CA
2	F	29	LYS	CA

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	Н	29	LYS	Mainchain
2	L	1	PHE	Mainchain

4.2 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	Α	163	0	149	1	0
1	С	163	0	149	2	0
1	Ε	163	0	149	3	0
1	G	163	0	149	2	0
1	I	170	0	154	1	0
1	K	163	0	149	2	0
2	В	244	0	234	5	0
2	D	254	0	241	6	0

Continued on next page...



Continued from previous page...

Mol	Chain	Non-H		H(added)	Clashes	Symm-Clashes
2	F	244	0	234	1	0
2	Н	244	0	234	4	0
2	J	244	0	233	4	0
2	L	244	0	234	4	0
3	A	7	0	6	0	0
3	С	7	0	6	0	0
3	E	7	0	6	0	0
3	G	7	0	6	0	0
3	I	7	0	6	0	0
3	K	7	0	6	0	0
4	J	1	0	0	0	0
4	L	1	0	0	0	0
5	J	1	0	0	0	0
5	L	1	0	0	0	0
6	A	4	0	0	0	0
6	В	9	0	0	1	0
6	С	9	0	0	0	0
6	D	10	0	0	0	0
6	Е	3	0	0	0	0
6	F	6	0	0	0	0
6	G	4	0	0	0	0
6	Н	10	0	0	0	0
6	I	7	0	0	1	0
6	J	13	0	0	0	0
6	K	8	0	0	0	0
6	L	8	0	0	0	0
All	All	2596	0	2345	27	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.

The worst 5 of 27 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$egin{array}{c} ext{Interatomic} \ ext{distance } (ext{Å}) \end{array}$	$egin{array}{c} ext{Clash} \ ext{overlap } (ext{Å}) \end{array}$
1:C:10:ILE:HD11	2:L:2:VAL:HA	1.66	0.76
1:K:14:TYR:CD1	1:K:14:TYR:N	2.58	0.69
2:B:25:PHE:HB3	2:D:25[A]:PHE:CE1	2.31	0.66
1:E:2:ILE:HD13	1:E:19:TYR:CE2	2.34	0.62
2:B:14:ALA:O	2:B:18:VAL:HG13	2.01	0.59

There are no symmetry-related clashes.



4.3 Torsion angles (i)

4.3.1 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	entiles
1	A	$19/21 \; (90\%)$	18 (95%)	1 (5%)	0	100	100
1	С	19/21 (90%)	19 (100%)	0	0	100	100
1	Е	19/21 (90%)	19 (100%)	0	0	100	100
1	G	19/21 (90%)	19 (100%)	0	0	100	100
1	I	20/21 (95%)	19 (95%)	1 (5%)	0	100	100
1	K	19/21 (90%)	18 (95%)	1 (5%)	0	100	100
2	В	28/30 (93%)	28 (100%)	0	0	100	100
2	D	30/30 (100%)	28 (93%)	0	2 (7%)	1	0
2	F	28/30 (93%)	27 (96%)	0	1 (4%)	3	2
2	Н	28/30 (93%)	27 (96%)	1 (4%)	0	100	100
2	J	$28/30 \ (93\%)$	27 (96%)	1 (4%)	0	100	100
2	L	28/30 (93%)	26 (93%)	2 (7%)	0	100	100
All	All	285/306~(93%)	275 (96%)	7 (2%)	3 (1%)	14	15

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	F	29	LYS
2	D	29	LYS
2	D	2	VAL

4.3.2 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	Perce	${f ntiles}$
1	A	$20/20 \; (100\%)$	18 (90%)	2 (10%)	7	9
1	С	$20/20 \; (100\%)$	18 (90%)	2 (10%)	7	9
1	E	$20/20 \; (100\%)$	18 (90%)	2 (10%)	7	9
1	G	$20/20 \; (100\%)$	17 (85%)	3 (15%)	3	3
1	I	$21/20 \ (105\%)$	19 (90%)	2 (10%)	8	10
1	K	$20/20 \; (100\%)$	18 (90%)	2 (10%)	7	9
2	В	$26/26 \; (100\%)$	24 (92%)	2 (8%)	13	16
2	D	$28/26 \ (108\%)$	24 (86%)	4 (14%)	3	3
2	F	$26/26 \; (100\%)$	22 (85%)	4 (15%)	2	2
2	Н	$26/26 \; (100\%)$	23 (88%)	3 (12%)	5	6
2	J	$26/26 \; (100\%)$	20 (77%)	6 (23%)	1	0
2	L	$26/26 \; (100\%)$	19 (73%)	7 (27%)	0	0
All	All	279/276 (101%)	240 (86%)	39 (14%)	3	3

5 of 39 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	G	15	GLN
2	Н	27	THR
2	L	22	ARG
1	G	21	ASN
2	Н	17	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	${f Res}$	\mathbf{Type}
1	A	5	GLN
1	С	5	GLN
1	G	15	GLN
2	J	3	ASN
1	K	21	ASN

4.3.3 RNA (i)

There are no RNA molecules in this entry.



4.4 Non-standard residues in protein, DNA, RNA chains (i)

There are no non-standard protein/DNA/RNA residues in this entry.

4.5 Carbohydrates (i)

There are no carbohydrates in this entry.

4.6 Ligand geometry (i)

Of 10 ligands modelled in this entry, 4 are monoatomic - leaving 6 for Mogul analysis.

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	T-m	Chain	Dog	Res Link Bond lengths			Bond angles			
MIOI	Type	Chain	m Res	Link	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	IPH	I	22	-	7,7,7	0.58	0	8,8,8	0.83	0
3	IPH	С	22	_	7,7,7	0.67	0	8,8,8	1.21	0
3	IPH	Е	22	_	7,7,7	0.75	0	8,8,8	1.08	1 (12%)
3	IPH	K	22	_	7,7,7	0.51	0	8,8,8	0.64	0
3	IPH	A	22	-	7,7,7	0.91	0	8,8,8	0.69	0
3	IPH	G	22	-	7,7,7	0.84	0	8,8,8	1.36	1 (12%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	IPH	I	22	_	_	_	0/1/1/1
3	IPH	С	22	_	-	-	0/1/1/1
3	IPH	Е	22	_	-	_	0/1/1/1
3	IPH	K	22	_	-	-	0/1/1/1
3	IPH	A	22	_	-	-	0/1/1/1
3	IPH	G	22	_	-	-	0/1/1/1

There are no bond length outliers.



All (2) bond angle outliers are listed below:

Mo	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^{o})$
3	E	22	IPH	C4-C3-C2	2.35	123.76	120.19
3	G	22	IPH	C3-C2-C1	-2.06	116.40	119.31

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

4.7 Other polymers (i)

There are no such residues in this entry.

4.8 Polymer linkage issues (i)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
2	J	1

All chain breaks are listed below:

Mod	el C	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1		J	28:PRO	С	29:LYS	N	1.14



5 Fit of model and data (i)

5.1 Protein, DNA and RNA chains (i)

EDS was not executed - this section is therefore empty.

5.2 Non-standard residues in protein, DNA, RNA chains (i)

EDS was not executed - this section is therefore empty.

5.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

5.4 Ligands (i)

EDS was not executed - this section is therefore empty.

5.5 Other polymers (i)

EDS was not executed - this section is therefore empty.

