

# wwPDB X-ray Structure Validation Summary Report (i)

### Aug 10, 2020 – 10:00 AM BST

PDB ID : 1HRN

Title : HIGH RESOLUTION CRYSTAL STRUCTURES OF RECOMBINANT HU-

MAN RENIN IN COMPLEX WITH POLYHYDROXYMONOAMIDE IN-

**HIBITORS** 

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Deposited on : 1995-03-31

Resolution : 1.80 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

buster-report : 1.1.7 (2018)

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)

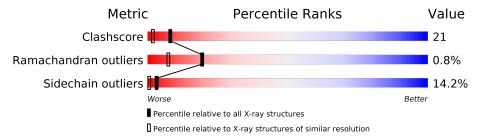
Validation Pipeline (wwPDB-VP) : 2.13.1

## 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 1.80 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	$(\# \mathrm{Entries})$	$(\#  ext{Entries},  ext{resolution range}( ext{Å}))$
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)

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.

Mol	Chain	Length	Quality o	f chain		
1	A	337	64%	24%	11%	<del>.</del>
1	В	337	53%	31%	13% •	



## 2 Entry composition (i)

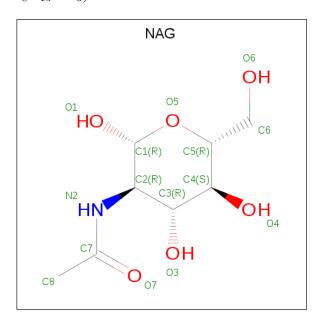
There are 4 unique types of molecules in this entry. The entry contains 5593 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 RENIN.

Mol	Chain	Residues	${f Atoms}$			ZeroOcc	AltConf	Trace		
1	Δ	335	Total C N O S		0	1				
1	11	333	2576	1644	417	501	14		U	1
1	D	334	Total	С	N	О	S	0	0	1
1	Б	354	2551	1630	409	498	14	0	0	1

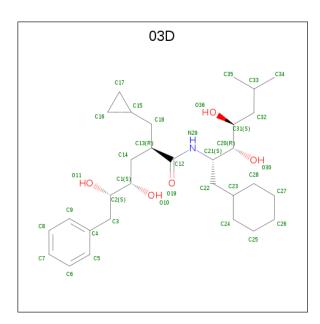
• Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total 14				0	0
2	В	1	Total 14	C 8	N 1	O 5	0	0

• Molecule 3 is (2R,4S,5S)-N-[(2S,3R,4S)-1-cyclohexyl-3,4-dihydroxy-6-methylheptan-2-yl]-2-(cyclopropylmethyl)-4,5-dihydroxy-6-phenylhexanamide (three-letter code: 03D) (formula:  $C_{30}H_{49}NO_5$ ).





Mol	Chain	Residues	${f Atoms}$				ZeroOcc	AltConf	
3	Δ	1	Total	С	N	О	0	0	
	Λ	1	36	30	1	5	U		
2	B	1	Total	С	Ν	Ο	0	0	
3	D	1	36	30	1	5	U	0	

### • Molecule 4 is water.

Mo	l	Chain	Residues	Atoms	ZeroOcc	AltConf
4		A	206	Total O 206 206	0	0
4		В	160	Total O 160 160	0	0

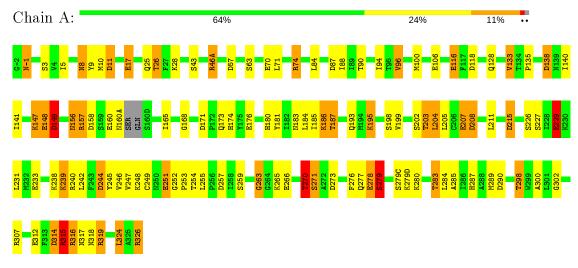


## 3 Residue-property plots (i)

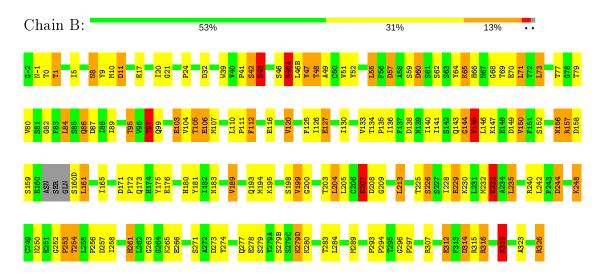
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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: RENIN



• Molecule 1: RENIN





# 4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants	141.10Å 141.10Å 141.10Å	Depositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	6.00 - 1.80	Depositor
% Data completeness	(Not available) (6.00-1.80)	Depositor
(in resolution range)	(1100 available) (0.00 1.00)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	TNT, X-PLOR	Depositor
$R, R_{free}$	0.215 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5593	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	27.0	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: NAG, 03D

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	Bond angles		
IVIOI	Chain	RMSZ	# Z  > 5	RMSZ	# Z >5	
1	A	1.17	15/2635~(0.6%)	1.51	38/3572 (1.1%)	
1	В	1.17	$14/2610 \ (0.5\%)$	1.57	40/3542 (1.1%)	
All	All	1.17	$29/5245 \ (0.6\%)$	1.54	78/7114 (1.1%)	

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 maintain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	В	1	0

The worst 5 of 29 bond length outliers are listed below:

Mol	Chain	${f Res}$	$\mathbf{Type}$	${f Atoms}$	$\mathbf{Z}$	$\operatorname{Observed}(\operatorname{\AA})$	$\mathbf{Ideal}(\mathbf{\AA})$
1	A	251	GLU	CD-OE1	11.52	1.38	1.25
1	A	233	GLU	CD-OE1	9.49	1.36	1.25
1	A	229	GLU	CD-OE1	9.06	1.35	1.25
1	В	17	GLU	CD-OE2	9.04	1.35	1.25
1	В	207	GLU	CD-OE1	8.75	1.35	1.25

The worst 5 of 78 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$\operatorname{Observed}(^o)$	$\mathbf{Ideal}(^{o})$
1	В	319	ARG	NE-CZ-NH1	17.15	128.87	120.30
1	В	319	ARG	NE-CZ-NH2	-15.55	112.53	120.30
1	В	11	ASP	CB-CG-OD1	11.18	128.36	118.30
1	A	46(A)	ARG	NE-CZ-NH1	10.88	125.74	120.30
1	В	314	ASP	CB-CG-OD2	-10.59	108.77	118.30



All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	В	95	THR	СВ

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	2576	0	2506	94	0
1	В	2551	0	2467	126	0
2	A	14	0	13	0	0
2	В	14	0	13	0	0
3	A	36	0	49	0	0
3	В	36	0	49	6	0
4	A	206	0	0	14	1
4	В	160	0	0	11	1
All	All	5593	0	5097	220	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 21.

The worst 5 of 220 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance}  ({\rm \AA}) \end{array}$	$egin{array}{c}  ext{Clash} \  ext{overlap } ( ext{Å}) \end{array}$
1:B:46(A):ARG:HG3	1:B:46(A):ARG:HH11	1.26	0.97
1:A:156:ASN:HD22	1:A:157:ARG:H	1.18	0.92
1:A:185:ILE:HG22	1:A:186:LYS:HG2	1.57	0.86
1:A:25:GLN:HE22	1:A:57:ASP:H	1.23	0.86
1:B:49:ALA:CB	1:B:107:MET:HG2	2.10	0.80

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	$egin{aligned}  ext{Interatomic} \  ext{distance} \ ( ext{Å}) \end{aligned}$	$egin{aligned}  ext{Clash} \  ext{overlap } ( ext{Å}) \end{aligned}$
4:A:927:HOH:O	4:B:917:HOH:O[6_456]	2.11	0.09

## 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	331/337 (98%)	315 (95%)	14 (4%)	2 (1%)	25 12
1	В	330/337~(98%)	315 (96%)	12 (4%)	3 (1%)	17 6
All	All	661/674 (98%)	630 (95%)	26 (4%)	5 (1%)	19 7

#### All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	279	SER
1	A	203	THR
1	В	242	LEU
1	В	112	PHE
1	В	253	PRO

#### 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	Analysed Rotameric Outliers		Percentiles			
1	A	$284/287 \ (99\%)$	244 (86%)	40 (14%)	3 0			
1	В	280/287~(98%)	240 (86%)	40 (14%)	3 0			
All	All	564/574 (98%)	484 (86%)	80 (14%)	3 0			



5 of 80 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	316	ARG
1	В	46	SER
1	В	235	LEU
1	A	319	ARG
1	В	1	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 21 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	318	ASN
1	В	99	GLN
1	В	261	HIS
1	A	317	ASN
1	В	277	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	Mol True C		Res	${ m Res} \mid { m Link} \mid$	Bo	ond leng	$ ag{ths}$	В	ond ang	les		
10101	Type	Chain	nes	nes	nes	Link	Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	03D	A	391	-	38,38,38	1.57	5 (13%)	41,51,51	1.18	5 (12%)		
2	NAG	A	367	1	14,14,15	0.95	1 (7%)	17,19,21	1.79	5 (29%)		
3	03D	В	391	-	38,38,38	1.52	5 (13%)	41,51,51	1.37	4 (9%)		
2	NAG	В	367	1	14,14,15	1.59	2 (14%)	17,19,21	1.98	7 (41%)		

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	03D	A	391	-	-	0/40/50/50	0/3/3/3
2	NAG	A	367	1	-	0/6/23/26	0/1/1/1
3	03D	В	391	-	-	1/40/50/50	0/3/3/3
2	NAG	В	367	1	-	3/6/23/26	0/1/1/1

The worst 5 of 13 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\operatorname{\AA})$	$\operatorname{Ideal}(\operatorname{\AA})$
3	A	391	03D	C20-C21	4.36	1.61	1.53
3	В	391	03D	C18-C15	4.00	1.59	1.53
3	A	391	03D	C22-C21	3.92	1.58	1.52
3	A	391	03D	C3-C4	3.60	1.60	1.51
2	В	367	NAG	C1-C2	3.38	1.57	1.52

The worst 5 of 21 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$\mathbf{Observed}(^o)$	$Ideal(^{o})$
3	В	391	03D	C14-C13-C12	-4.17	102.07	109.44
2	A	367	NAG	C6-C5-C4	-3.72	104.29	113.00
2	В	367	NAG	O4-C4-C3	3.57	118.60	110.35
3	В	391	03D	O19-C12-C13	-3.46	117.61	122.12
3	В	391	03D	C4-C3-C2	-3.45	107.44	113.61

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	В	367	NAG	C8-C7-N2-C2
2	В	367	NAG	O7-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
2	В	367	NAG	O5-C5-C6-O6
3	В	391	03D	C12-C13-C18-C15

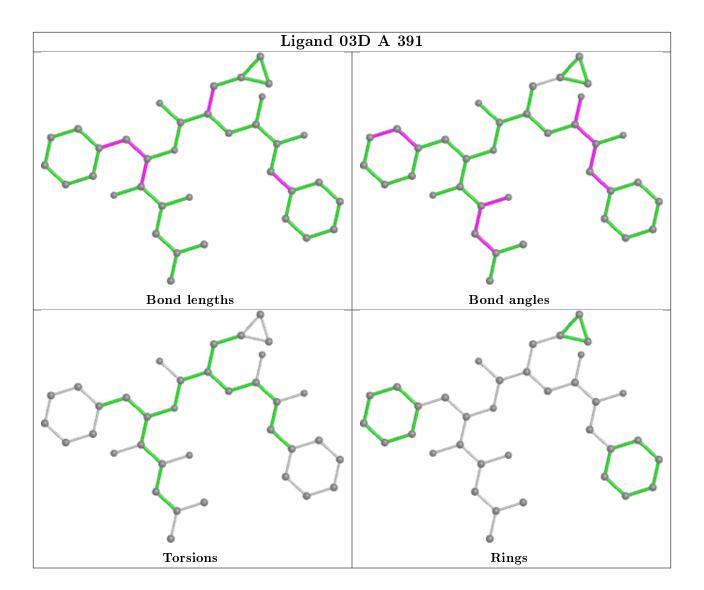
There are no ring outliers.

1 monomer is involved in 6 short contacts:

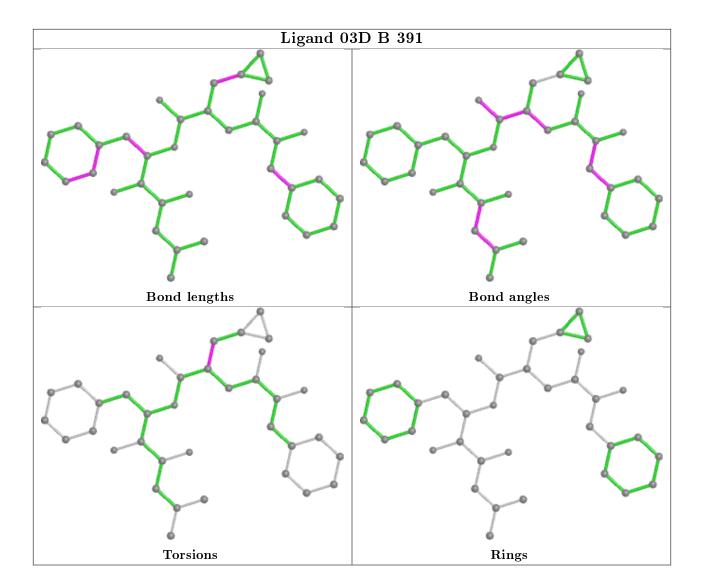
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	В	391	03D	6	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.









# 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)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

### 6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

## 6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

### 6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.

