

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

#### Oct 17, 2023 – 04:17 AM EDT

PDB ID : 1X2T

Title : Crystal Structure of Habu IX-bp at pH 6.5

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Deposited on : 2005-04-26

Resolution : 1.72 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 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 : 2.36

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

Refmac : 5.8.0158

CCP4 : 7.0.044 (Gargrove) roteins) : Engh & Huber (2001)

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

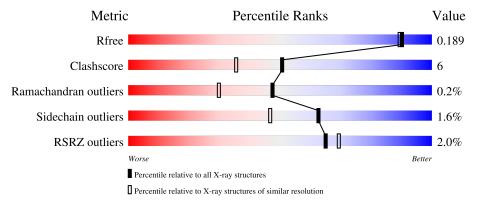
Validation Pipeline (wwPDB-VP) : 2.36

## 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 1.72 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	$(\#  ext{Entries})$	$(\#  ext{Entries},  ext{ resolution range}( ext{Å}))$
$R_{free}$	130704	5722 (1.74-1.70)
Clashscore	141614	6152 (1.74-1.70)
Ramachandran outliers	138981	6051 (1.74-1.70)
Sidechain outliers	138945	6051 (1.74-1.70)
RSRZ outliers	127900	5629 (1.74-1.70)

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	A	129	84%	16%	•
1	С	129	84%	16%	-
2	В	123	84%	16%	_
2	D	123	87%	13%	_



## 2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 4698 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 Coagulation factor IX/X-binding protein A chain.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Δ	129	Total	С	N	О	S	0	1	0
1	11	125	1033	649	170	207	7		1	0
1	C	129	Total	С	N	О	S	0	1	0
1		129	1033	649	170	207	7	0	1	U

• Molecule 2 is a protein called Coagulation factor IX/factor X-binding protein B chain.

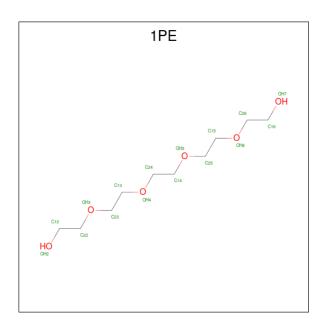
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace	
2	В	123	Total 1017		N 173	S 11	0	0	0
2	D	123	Total 1017		N 173	S 11	0	0	0

• Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Ca 1 1	0	0
3	В	1	Total Ca 1 1	0	0
3	С	1	Total Ca 1 1	0	0
3	D	1	Total Ca 1 1	0	0

• Molecule 4 is PENTAETHYLENE GLYCOL (three-letter code: 1PE) (formula: C<sub>10</sub>H<sub>22</sub>O<sub>6</sub>).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	В	1	Total C O 13 8 5	0	0
4	D	1	Total C O 13 8 5	0	0

### • Molecule 5 is water.

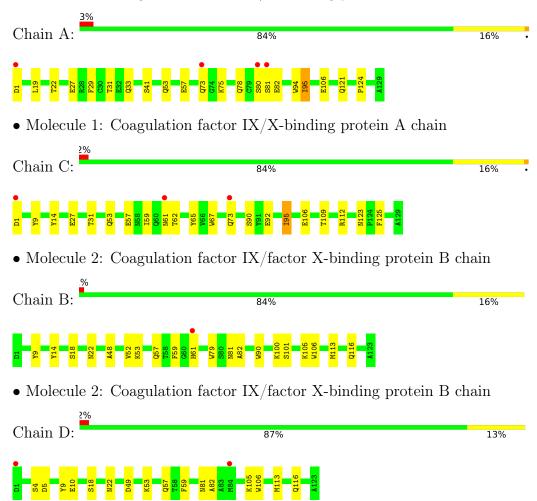
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	149	Total O 149 149	0	0
5	В	136	Total O 136 136	0	0
5	С	122	Total O 122 122	0	0
5	D	161	Total O 161 161	0	0



## 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: Coagulation factor IX/X-binding protein A chain





# 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants	60.71Å 63.51Å 66.90Å	Donogiton
a, b, c, $\alpha$ , $\beta$ , $\gamma$	90.00° 117.00° 90.00°	Depositor
Resolution (Å)	33.26 - 1.72	Depositor
Resolution (A)	54.11 - 1.72	EDS
% Data completeness	99.0 (33.26-1.72)	Depositor
(in resolution range)	99.0 (54.11-1.72)	EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$< I/\sigma(I) > 1$	2.43 (at 1.72Å)	Xtriage
Refinement program	CNS 1.1	Depositor
$R, R_{free}$	0.191 , 0.222	Depositor
$1\iota$ , $1\iota_{free}$	0.184 , $0.189$	DCC
$R_{free}$ test set	2421 reflections $(5.06\%)$	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	15.2	Xtriage
Anisotropy	0.544	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.32 , 44.4	EDS
L-test for twinning <sup>2</sup>	$< L > = 0.49, < L^2> = 0.32$	Xtriage
Estimated twinning fraction	0.003 for h,-k,-h-l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	4698	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 18.86% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;sup>2</sup>Theoretical values of <|L|>,  $<L^2>$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



<sup>&</sup>lt;sup>1</sup>Intensities 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: 1PE, CA

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	ol Chain	Bond	lengths	Bond angles		
MIOI		RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	A	0.33	0/1066	0.63	1/1440 (0.1%)	
1	С	0.31	0/1066	0.59	1/1440 (0.1%)	
2	В	0.34	0/1054	0.53	0/1427	
2	D	0.34	0/1054	0.54	0/1427	
All	All	0.33	0/4240	0.57	$2/5734 \ (0.0\%)$	

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	$\operatorname{Res}$	Type	Atoms	${f Z}$	$\mathbf{Observed}(^{o})$	$\mathbf{Ideal}(^o)$
1	A	95	ILE	N-CA-C	-6.59	93.19	111.00
1	С	95	ILE	N-CA-C	-5.73	95.52	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	A	1033	0	948	11	0
1	С	1033	0	948	15	0
2	В	1017	0	905	13	0
2	D	1017	0	905	11	0
3	A	1	0	0	0	0

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Continued	trom	mmoninonic	maaa
COHABABACA		DIEUIUU	DUIUE
0 0 1000100000			

Mol	Chain	Non-H	H(model)	$\mathbf{H}(\mathbf{added})$	Clashes	Symm-Clashes
3	В	1	0	0	0	0
3	С	1	0	0	0	0
3	D	1	0	0	0	0
4	В	13	0	17	0	0
4	D	13	0	17	0	0
5	A	149	0	0	1	0
5	В	136	0	0	2	0
5	С	122	0	0	2	0
5	D	161	0	0	2	0
All	All	4698	0	3740	45	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 45 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	Clash overlap (Å)
2:B:18:SER:HA	2:B:116:GLN:HE21	1.32	0.95
2:D:18:SER:HA	2:D:116:GLN:HE21	1.36	0.91
1:C:95:ILE:HG13	2:D:105:LYS:HB3	1.76	0.67
2:B:18:SER:CA	2:B:116:GLN:HE21	2.08	0.64
2:B:18:SER:HA	2:B:116:GLN:NE2	2.09	0.62

There are no symmetry-related clashes.

## 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	A	128/129 (99%)	124 (97%)	4 (3%)	0	100	100
1	С	128/129 (99%)	123 (96%)	5 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
2	В	121/123 (98%)	115 (95%)	5 (4%)	1 (1%)	19 6
2	D	121/123 (98%)	114 (94%)	7 (6%)	0	100 100
All	All	498/504 (99%)	476 (96%)	21 (4%)	1 (0%)	47 30

#### All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	В	101	SER

#### 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	112/111 (101%)	108 (96%)	4 (4%)	35 15
1	$\mathbf{C}$	112/111 (101%)	111 (99%)	1 (1%)	78 69
2	В	106/106 (100%)	104 (98%)	2 (2%)	57 39
2	D	106/106 (100%)	106 (100%)	0	100 100
All	All	436/434 (100%)	429 (98%)	7 (2%)	62 47

5 of 7 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	106	GLU
2	В	90	TRP
1	С	1	ASP
2	В	100	LYS
1	A	78	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 10 such sidechains are listed below:

Mol	Chain	Res	Type
1	С	123	ASN

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Mol	Chain	Res	Type
2	D	32	GLN
2	D	57	GLN
2	В	78	GLN
2	В	116	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)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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	Type	Chain	Chain	Chain	Res	Link	Bond lengths			Bond angles		
WIOI	туре	Chain	nes	LIIIK	Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2		
4	1PE	В	602	-	12,12,15	0.92	0	11,11,14	2.23	4 (36%)		
4	1PE	D	601	-	12,12,15	0.93	0	11,11,14	2.23	4 (36%)		

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
Г	4	1PE	В	602	-	-	1/10/10/13	-
	4	1PE	D	601	-	-	3/10/10/13	-

There are no bond length outliers.

The worst 5 of 8 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$\mathbf{Observed}(^o)$	$Ideal(^{o})$
4	В	602	1PE	OH4-C13-C23	-4.45	90.52	110.07
4	D	601	1PE	OH4-C13-C23	-4.21	91.56	110.07
4	D	601	1PE	OH6-C15-C25	-3.59	94.22	110.39
4	D	601	1PE	C24-OH4-C13	-3.44	98.39	113.29
4	В	602	1PE	OH6-C15-C25	-3.33	95.38	110.39

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	В	602	1PE	OH5-C14-C24-OH4
4	D	601	1PE	ОН7-С16-С26-ОН6
4	D	601	1PE	OH5-C14-C24-OH4
4	D	601	1PE	ОН6-С15-С25-ОН5

There are no ring outliers.

No monomer is involved in short contacts.

## 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	$\langle { m RSRZ} \rangle$	$\# \mathrm{RSRZ}{>}2$	$OWAB(Å^2)$	Q<0.9
1	A	129/129 (100%)	0.02	4 (3%) 49 53	7, 13, 30, 40	0
1	С	129/129 (100%)	0.02	3 (2%) 60 65	10, 17, 29, 37	0
2	В	123/123 (100%)	-0.01	1 (0%) 86 89	9, 16, 25, 31	0
2	D	123/123 (100%)	-0.08	2 (1%) 72 76	9, 15, 23, 34	0
All	All	504/504 (100%)	-0.01	10 (1%) 65 69	7, 15, 27, 40	0

The worst 5 of 10 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	81	SER	5.7
1	С	61	ASN	3.5
1	A	73	GLN	3.3
2	В	61	HIS	3.0
2	D	84	MET	2.9

## 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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
4	1PE	В	602	13/16	0.78	0.20	49,52,56,56	0
4	1PE	D	601	13/16	0.86	0.18	16,33,47,49	0
3	CA	В	604	1/1	0.98	0.06	14,14,14,14	0
3	CA	С	605	1/1	0.98	0.06	17,17,17,17	0
3	CA	A	603	1/1	0.99	0.06	19,19,19,19	0
3	CA	D	606	1/1	1.00	0.06	10,10,10,10	0

# 6.5 Other polymers (i)

There are no such residues in this entry.

