

wwPDB X-ray Structure Validation Summary Report (i)

May 24, 2020 – 11:05 pm BST

PDB ID : 1JRR

Title: HUMAN PLASMINOGEN ACTIVATOR INHIBITOR-2.[LOOP (66-98)

DELETIONMUTANT COMPLEXED WITH PEPTIDE MIMICKING THE

REACTIVE CENTER LOOP

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Deposited on : 2001-08-15

Resolution : 1.60 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)

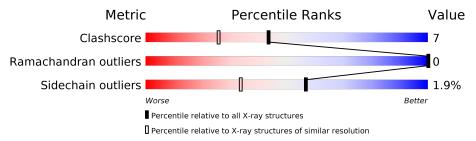
Validation Pipeline (wwPDB-VP) : 2.11

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 1.60 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	3665 (1.60-1.60)		
Ramachandran outliers	138981	3564 (1.60-1.60)		
Sidechain outliers	138945	3563 (1.60-1.60)		

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 of chain							
1	A	382	81%	12%	• 7%					
2	Р	15	100%							



2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 3419 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 PLASMINOGEN ACTIVATOR INHIBITOR-2.

Mol	Chain	Residues	${f Atoms}$			ZeroOcc	AltConf	Trace		
1	Λ	357	Total	С	N	О	S	0	19	0
1	A	397	2886	1840	476	551	19	0	10	

There are 33 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	ASN	DELETION	UNP P05120
A	?	-	ALA	DELETION	UNP P05120
A	?	-	VAL	DELETION	UNP P05120
A	?	-	THR	DELETION	UNP P05120
A	?	-	PRO	DELETION	UNP P05120
A	?	-	MET	DELETION	UNP P05120
A	?	-	THR	DELETION	UNP P05120
A	?	-	PRO	DELETION	UNP P05120
A	?	-	GLU	DELETION	UNP P05120
A	?	-	ASN	DELETION	UNP P05120
A	?	-	PHE	DELETION	UNP P05120
A	?	-	THR	DELETION	UNP P05120
A	?	-	SER	DELETION	UNP P05120
A	?	-	CYS	DELETION	UNP P05120
A	?	-	GLY	DELETION	UNP P05120
A	?	-	PHE	DELETION	UNP P05120
A	?	-	MET	DELETION	UNP P05120
A	?	-	GLN	DELETION	UNP P05120
A	?	-	GLN	DELETION	UNP P05120
A	?	-	ILE	DELETION	UNP P05120
A	?	-	GLN	DELETION	UNP P05120
A	?	-	LYS	DELETION	UNP P05120
A	?	-	GLY	DELETION	UNP P05120
A	?	-	SER	DELETION	UNP P05120
A	?	-	TYR	DELETION	UNP P05120
A	?	-	PRO	DELETION	UNP P05120
A	?	-	ASP	DELETION	UNP P05120

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Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	ALA	DELETION	UNP P05120
A	?	-	ILE	DELETION	UNP P05120
A	?	-	LEU	DELETION	UNP P05120
A	?	-	GLN		UNP P05120
A	?	-	ALA	DELETION	UNP P05120
A	?	-	GLN	DELETION	UNP P05120

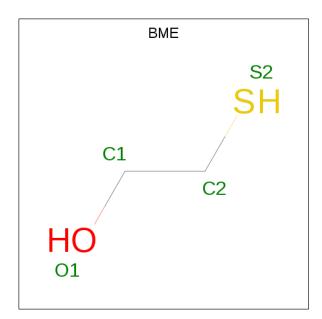
• Molecule 2 is a protein called PLASMINOGEN ACTIVATOR INHIBITOR-2.

Mol	Chain	Residues	${f Atoms}$			ZeroOcc	AltConf	Trace		
9	D	15	Total	С	N	О	S	0	0	0
	Γ	19	91	52	17	21	1	0	0	U

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference	
Р	366	ACE	-	ACETYLATION	UNP P05120	

• Molecule 3 is BETA-MERCAPTOETHANOL (three-letter code: BME) (formula: C₂H₆OS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total 4	C 2	O 1	S 1	0	0

• Molecule 4 is water.



I	Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
	4	A	433	Total O 433 433	0	0
	4	Р	5	Total O 5 5	0	0

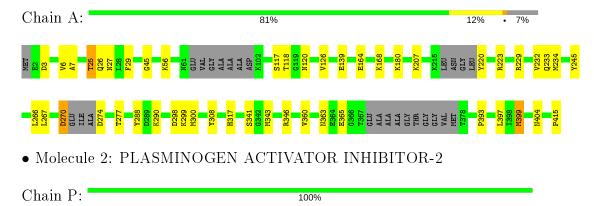


3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA 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: PLASMINOGEN ACTIVATOR INHIBITOR-2



There are no outlier residues recorded for this chain.



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 21 2	Depositor	
Cell constants	93.14Å 104.16Å 41.40Å	Depositor	
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor	
Resolution (Å)	69.01 - 1.60	Depositor	
% Data completeness	94.6 (69.01-1.60)	Depositor	
(in resolution range)	`	Depositor	
R_{merge}	(Not available)	Depositor	
R_{sym}	0.05	Depositor	
Refinement program	REFMAC 5.0	Depositor	
R, R_{free}	0.197 , 0.225	Depositor	
Estimated twinning fraction	No twinning to report.	Xtriage	
Total number of atoms	3419	wwPDB-VP	
Average B, all atoms (Å ²)	21.0	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: ACE, BME

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		
MIOI		RMSZ	# Z > 5	RMSZ	# Z > 5	
1	A	0.59	0/3012	0.87	7/4052 (0.2%)	
2	Р	0.58	0/88	0.80	0/116	
All	All	0.59	0/3100	0.86	7/4168 (0.2%)	

There are no bond length outliers.

The worst 5 of 7 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^o)$	$\mathbf{Ideal}(^{o})$
1	Α	139[A]	GLU	CA-CB-CG	6.66	128.06	113.40
1	Α	139[B]	GLU	CA-CB-CG	6.66	128.06	113.40
1	A	274	ASP	CB-CG-OD2	6.37	124.04	118.30
1	A	126[A]	VAL	CA-CB-CG2	6.00	119.90	110.90
1	A	126[B]	VAL	CA-CB-CG2	6.00	119.90	110.90

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.

\mathbf{N}	Iol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
	1	Α	2886	0	2849	40	0
	2	Р	91	0	88	0	0
	3	A	4	0	5	0	0
	4	A	433	0	0	12	4

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Mol	Chain	Non-H	H(model)	$\mathbf{H}(\mathbf{added})$	Clashes	Symm-Clashes
4	Р	5	0	0	0	0
All	All	3419	0	2942	40	4

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 7.

The worst 5 of 40 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$egin{array}{l} ext{Interatomic} \ ext{distance} \ (ext{Å}) \end{array}$	Clash overlap (Å)
1:A:346:ARG:HD2	4:A:830:HOH:O	1.81	0.81
1:A:26:GLN:HE21	1:A:27:ASN:H	1.35	0.73
1:A:397:LEU:HD13	1:A:399[A]:MET:SD	2.36	0.66
1:A:168:LYS:HE3	4:A:987:HOH:O	1.98	0.64
1:A:207:LYS:HE3	1:A:365:GLU:HG2	1.83	0.59

All (4) 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	$egin{array}{l} ext{Interatomic} \ ext{distance} \ (ext{Å}) \end{array}$	$egin{aligned} ext{Clash} \ ext{overlap } (ext{Å}) \end{aligned}$
4:A:658:HOH:O	4:A:659:HOH:O[2_655]	2.03	0.17
4:A:705:HOH:O	4:A:824:HOH:O[3_545]	2.04	0.16
4:A:704:HOH:O	4:A:823:HOH:O[3_545]	2.15	0.05
4:A:815:HOH:O	4:A:816:HOH:O[3_555]	2.19	0.01

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	\mathbf{ntiles}
1	A	$360/382 \ (94\%)$	350 (97%)	10 (3%)	0	100	100
2	Р	13/15 (87%)	13 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed			
All	All	373/397 (94%)	363 (97%)	10 (3%)	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	$329/329 \ (100\%)$	320 (97%)	9 (3%)	44 20		
2	Р	7/7 (100%)	7 (100%)	0	100 100		
All	All	336/336 (100%)	327 (97%)	9 (3%)	57 20		

5 of 9 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	317	HIS
1	A	399[B]	MET
1	A	341[B]	SER
1	A	117[B]	SER
1	A	341[A]	SER

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

Mol	Chain	${f Res}$	\mathbf{Type}
1	A	14	ASN
1	A	26	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 carbohydrates in this entry.

5.6 Ligand geometry (i)

1 ligand is 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	B	ond leng	$_{ m gths}$	Е	ond ang	gles
WIOI	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	BME	A	561	1	3,3,3	0.27	0	1,2,2	0.18	0

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	${ m Res}$	Link	Chirals	Torsions	Rings
3	BME	A	561	1	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	561	BME	O1-C1-C2-S2

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)

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.

