

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

#### May 29, 2020 - 12:51 am BST

PDB ID	:	2PTN
$\operatorname{Title}$	:	ON THE DISORDERED ACTIVATION DOMAIN IN TRYPSINOGEN.
		CHEMICAL LABELLING AND LOW-TEMPERATURE CRYSTALLOG-
		RAPHY
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Deposited on		
Resolution	:	1.55  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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

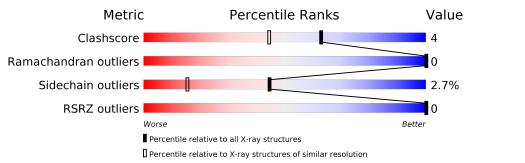
$\operatorname{MolProbity}$	:	4.02b-467
Xtriage (Phenix)	:	1.13
$\mathrm{EDS}$	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
$\operatorname{Refmac}$	:	5.8.0158
$\operatorname{CCP4}$	:	$7.0.044 (\mathrm{Gargrove})$
Ideal geometry (proteins)	:	Engh & 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 1.55 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
Clashscore	141614	1529 (1.56-1.56)
Ramachandran outliers	138981	1498 (1.56-1.56)
Sidechain outliers	138945	1495(1.56-1.56)
RSRZ outliers	127900	1465(1.56-1.56)

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% 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	А	223	86%	13%	•



# 2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 1712 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 TRYPSIN.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
1	А	223	Total	С	Ν	Ο	$\mathbf{S}$	65	0	0
1	11	220	1629	1012	279	324	14	00	0	0

• Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	А	1	Total Ca 1 1	0	0

• Molecule 3 is water.

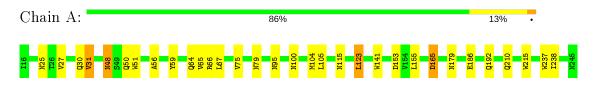
Mol	Chain	Residues Atoms		ZeroOcc	AltConf
3	А	82	Total         O           82         82	0	0



# 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 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: TRYPSIN





## 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	54.89Å $58.52$ Å $67.63$ Å	Depositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $90.00^{\circ}$ $90.00^{\circ}$	Depositor
Resolution (Å)	6.50 - 1.55	Depositor
Resolution (A)	27.45 - 1.50	EDS
% Data completeness	(Not available) $(6.50-1.55)$	Depositor
(in resolution range)	67.2(27.45-1.50)	EDS
R <sub>merge</sub>	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$< I/\sigma(I) >$	-	Xtriage
Refinement program	unknown	Depositor
$R, R_{free}$	0.193 , (Not available)	Depositor
It, It <sub>free</sub>	0.190 , (Not available)	DCC
$R_{free}$ test set	No test flags present.	wwPDB-VP
Wilson B-factor ( $Å^2$ )	10.5	Xtriage
Anisotropy	0.537	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.30 , $52.0$	EDS
L-test for twinning <sup>1</sup>	$ \langle L  \rangle = 0.41, \langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	1712	wwPDB-VP
Average B, all atoms $(Å^2)$	14.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 9.54% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;sup>1</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



# 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: 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	Chain	Bo	nd lengths	Bond angles		
		RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	А	1.01	5/1660~(0.3%)	0.96	3/2250~(0.1%)	

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

Mol	Chain   #Chirality outliers   #Planarity o		#Planarity outliers
1	А	0	8

Chain  $\mathbf{Z}$ Observed(Å) Mol  $\mathbf{Res}$ Type Atoms Ideal(Å) 1 А 215TRP NE1-CE2 -7.661.271.371 А 141 $\mathrm{TRP}$ NE1-CE2 -7.551.271.371 А 51TRP NE1-CE2 -7.511.271.37237 TRP NE1-CE2 1 А -7.371.271.37GLU 1 А 186CD-OE25.881.321.25

All (5) bond length outliers are listed below:

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^{o})$	$\mathbf{Ideal}(^{o})$
1	А	75	VAL	CA-CB-CG2	5.95	119.82	110.90
1	А	31	VAL	CA-CB-CG1	5.23	118.74	110.90
1	А	165	ASP	CB-CG-OD2	-5.09	113.72	118.30

There are no chirality outliers.

All (8) planarity outliers are listed below:



Mol	Chain	Res	Type	Group
1	А	115	ASN	Sidechain
1	А	165	ASP	Sidechain
1	А	210	GLN	Sidechain
1	А	25	ASN	Sidechain
1	А	27	VAL	Mainchain
1	А	30	GLN	Sidechain
1	А	79	ASN	Sidechain
1	А	95	ASN	Sidechain

### 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	А	1629	0	1588	13	1
2	А	1	0	0	0	0
3	А	82	0	0	0	0
All	All	1712	0	1588	13	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 4.

All (13) 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:48:ASN:HD22	1:A:50:GLN:H	1.41	0.69
1:A:31:VAL:HG22	1:A:67:LEU:CD2	2.34	0.57
1:A:31:VAL:HG13	1:A:65:VAL:HG13	1.94	0.49
1:A:31:VAL:HG22	1:A:67:LEU:HD23	1.93	0.49
1:A:48:ASN:ND2	1:A:50:GLN:H	2.11	0.49
1:A:64:GLN:HE22	1:A:66:ARG:HH21	1.63	0.47
1:A:123:LEU:HD12	1:A:123:LEU:HA	1.86	0.45
1:A:123:LEU:HD21	1:A:238:ILE:HG21	2.01	0.43
1:A:48:ASN:HD22	1:A:48:ASN:C	2.23	0.42
1:A:56:ALA:HA	1:A:104:MET:HB2	2.02	0.41
1:A:100:ASN:HD21	1:A:179:ASN:HD22	1.68	0.41
1:A:64:GLN:HE21	1:A:66:ARG:HE	1.68	0.41
1:A:64:GLN:NE2	1:A:66:ARG:HE	2.19	0.40



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	Interatomic distance (Å)	Clash overlap (Å)
1:A:59:TYR:OH	$1:A:153:ASP:OD2[4_556]$	2.17	0.03

### 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	А	221/223~(99%)	217~(98%)	4 (2%)	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	А	184/184~(100%)	179~(97%)	5(3%)	44 15		

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	$\mathbf{Res}$	Type
1	А	48	ASN
1	А	105	LEU
1	А	123	LEU
1	А	155	LEU
1	А	192	GLN



Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	А	30	GLN
1	А	48	ASN
1	А	64	GLN
1	А	97	ASN
1	А	100	ASN
1	А	101	ASN
1	А	210	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)

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

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<sup>th</sup> 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	<RSRZ $>$	#	₽RSF	RZ>2	$OWAB(Å^2)$	Q<0.9
1	А	223/223~(100%)	-0.17	0	100	100	7, 13, 24, 34	38 (17%)

There are no RSRZ outliers to report.

### 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 carbohydrates 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	${f B} ext{-factors}({ m \AA}^2)$	$Q{<}0.9$
2	CA	А	480	1/1	0.99	0.04	$11,\!11,\!11,\!11$	0

### 6.5 Other polymers (i)

There are no such residues in this entry.

