



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 31, 2020 – 09:02 AM BST

PDB ID : 1BZX
Title : THE CRYSTAL STRUCTURE OF ANIONIC SALMON TRYPSIN IN COM-
PLEX WITH BOVINE PANCREATIC TRYPSIN INHIBITOR
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Deposited on : 1998-11-05
Resolution : 2.10 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 ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	FAILED
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

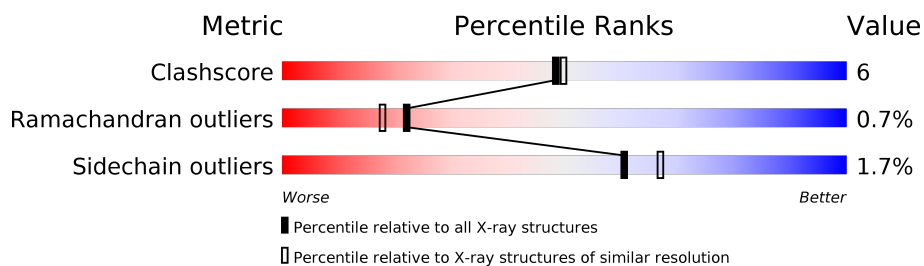
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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 (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 2219 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 PROTEIN (TRYPSIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	E	222	Total	C	N	O	S	58	3	0
			1675	1046	280	330	19			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	24	PRO	ALA	CONFLICT	UNP P35031
E	28	PRO	THR	CONFLICT	UNP P35031

- Molecule 2 is a protein called PROTEIN (BOVINE PANCREATIC TRYPSIN INHIBITOR).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	I	58	Total	C	N	O	S	19	1	0
			458	287	84	79	8			

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	1	Total	Ca	0	0
			1	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	E	66	Total	O	0	0
			66	66		
4	I	19	Total	O	0	0
			19	19		

SEQUENCE-PLOTS INFOmissingINFO

3 Data and refinement statistics

EDS failed to run properly - this section is therefore incomplete.

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	84.12Å 84.12Å 222.15Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	8.00 – 2.10	Depositor
% Data completeness (in resolution range)	97.3 (8.00-2.10)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.59 (at 2.14Å)	Xtriage
Refinement program	X-PLOR 3.8	Depositor
R, R_{free}	0.206 , 0.238	Depositor
Wilson B-factor (Å ²)	38.3	Xtriage
Anisotropy	0.025	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2219	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.42% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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.

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: 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	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	E	0.49	0/1728	0.70	0/2351
2	I	0.59	1/473 (0.2%)	0.71	0/632
All	All	0.52	1/2201 (0.0%)	0.71	0/2983

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	I	515	LYS	C-N	-6.07	1.20	1.34

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	E	1675	0	1582	20	0
2	I	458	0	439	5	0
3	E	1	0	0	0	0
4	E	66	0	0	1	0
4	I	19	0	0	0	0
All	All	2219	0	2021	24	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.

All (24) 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:E:74:LYS:HE3	1:E:74:LYS:HA	1.69	0.74
2:I:504:PHE:O	2:I:507:GLU:HG2	2.00	0.62
2:I:502:PRO:HG2	2:I:505:CYS:SG	2.40	0.60
1:E:169:ASN:HD21	1:E:174:GLY:H	1.53	0.55
2:I:535:TYR:CZ	2:I:540:ALA:HB2	2.42	0.55
1:E:114:LEU:HA	1:E:118:VAL:O	2.11	0.49
1:E:97:TYR:O	2:I:539:ARG:NH2	2.46	0.49
1:E:90:ARG:HA	1:E:104:MET:HB2	1.95	0.49
1:E:210:GLN:NE2	4:E:781:HOH:O	2.46	0.49
1:E:26:SER:C	1:E:28:PRO:HD3	2.32	0.48
1:E:32:SER:OG	1:E:40:HIS:HD2	1.97	0.47
1:E:113:THR:O	1:E:115:ASN:ND2	2.48	0.47
1:E:60:LYS:HB2	1:E:63[B]:VAL:CG1	2.47	0.45
1:E:129:ALA:HA	1:E:130:PRO:HD3	1.67	0.45
1:E:74:LYS:HE3	1:E:74:LYS:CA	2.43	0.44
1:E:67:LEU:HD12	1:E:81:GLN:O	2.18	0.44
1:E:91:HIS:HA	1:E:237:TRP:CZ2	2.53	0.43
2:I:520:ARG:O	2:I:532:THR:HA	2.18	0.43
1:E:158:LEU:HD11	1:E:188:LYS:HB3	2.01	0.43
1:E:16:ILE:O	1:E:144:THR:HA	2.19	0.42
1:E:72:ASN:HA	1:E:153:ASN:O	2.19	0.42
1:E:40:HIS:HE1	1:E:193:GLY:O	2.02	0.42
1:E:199:VAL:HG21	1:E:228:TYR:CD2	2.54	0.42
1:E:169:ASN:ND2	1:E:174:GLY:H	2.16	0.40

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	Percentiles	
1	E	223/222 (100%)	215 (96%)	7 (3%)	1 (0%)	34	32
2	I	57/58 (98%)	56 (98%)	0	1 (2%)	8	4
All	All	280/280 (100%)	271 (97%)	7 (2%)	2 (1%)	22	18

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	49	GLU
2	I	556	GLY

4.3.2 Protein sidechains ⓘ

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	E	190/187 (102%)	186 (98%)	4 (2%)	53	59
2	I	47/46 (102%)	47 (100%)	0	100	100
All	All	237/233 (102%)	233 (98%)	4 (2%)	60	67

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	37	SER
1	E	61	SER
1	E	74	LYS
1	E	233	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	30	GLN
1	E	40	HIS
1	E	169	ASN
1	E	192	GLN
1	E	202	ASN

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Mol	Chain	Res	Type
1	E	210	GLN
1	E	224	ASN
2	I	531	GLN

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 monosaccharides in this entry.

4.6 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.

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	I	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	I	515:LYS	C	516:ALA	N	1.20

5 Fit of model and data

5.1 Protein, DNA and RNA chains

EDS failed to run properly - this section is therefore empty.

5.2 Non-standard residues in protein, DNA, RNA chains

EDS failed to run properly - this section is therefore empty.

5.3 Carbohydrates

EDS failed to run properly - this section is therefore empty.

5.4 Ligands

EDS failed to run properly - this section is therefore empty.

5.5 Other polymers

EDS failed to run properly - this section is therefore empty.