



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 3MFJ / pdb_00003mfj
Title : Bovine trypsin at 0.8 Å resolution, restrained refinement
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Deposited on : 2010-04-02
Resolution : 0.80 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

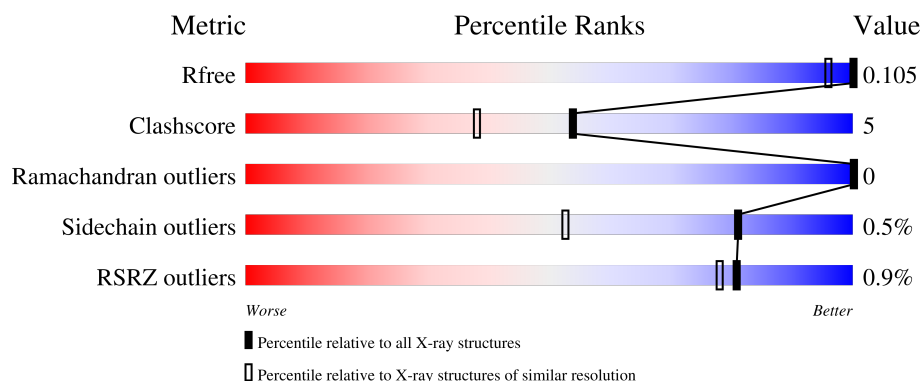
The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

i

X-RAY DIFFRACTION

A.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1419 (1.00-0.60)
Clashscore	190562	1475 (1.00-0.60)
Ramachandran outliers	187476	1389 (1.00-0.60)
Sidechain outliers	187428	1390 (1.00-0.60)
RSRZ outliers	180081	1412 (1.00-0.60)

Mol	Chain	Length	Quality of chain
1	A	223	89% 10%

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	BEN	A	252	-	X	-	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 3660 atoms, of which 1434 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 Cationic trypsin.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	223	Total	C	H	N	O	S	0	27	0
			3159	1074	1434	287	347	17			

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

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

- Molecule 3 is BENZAMIDINE (CCD ID: BEN) (formula: C₇H₈N₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	N	0	0
			9	7	2		

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	3	3		

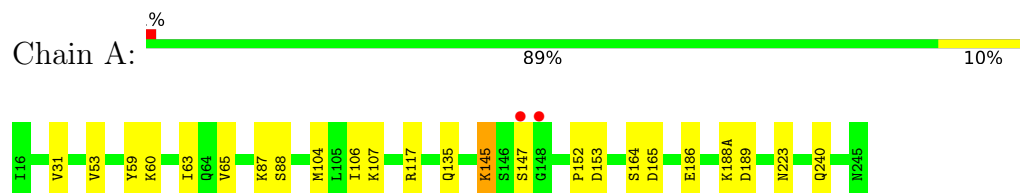
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	480	Total 480	O 480	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: Cationic trypsin



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	54.22Å 58.53Å 66.45Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 0.80 30.00 – 0.80	Depositor EDS
% Data completeness (in resolution range)	80.3 (30.00-0.80) 93.4 (30.00-0.80)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.38 (at 0.80Å)	Xtriage
Refinement program	SHELXL-97	Depositor
R, R_{free}	0.105 , 0.112 (Not available) , 0.105	Depositor DCC
R_{free} test set	2204 reflections (1.00%)	wwPDB-VP
Wilson B-factor (Å ²)	7.3	Xtriage
Anisotropy	0.117	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.48 , 126.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.99	EDS
Total number of atoms	3660	wwPDB-VP
Average B, all atoms (Å ²)	11.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.46% 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.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, BEN, CA, SO4

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	A	0.91	3/1867 (0.2%)	1.08	13/2529 (0.5%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	53[A]	VAL	CA-CB	-6.61	1.46	1.54
1	A	53[B]	VAL	CA-CB	-6.61	1.46	1.54
1	A	147	SER	C-O	-5.57	1.17	1.24

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	147	SER	CA-C-N	-7.57	108.92	121.57
1	A	147	SER	C-N-CA	-7.57	108.92	121.57
1	A	189	ASP	CA-CB-CG	6.55	119.15	112.60
1	A	117	ARG	NE-CZ-NH2	-6.17	113.64	119.20
1	A	135[A]	GLN	OE1-CD-NE2	-6.05	116.55	122.60
1	A	135[B]	GLN	OE1-CD-NE2	-6.05	116.55	122.60
1	A	145	LYS	CB-CG-CD	5.83	124.72	111.30
1	A	223	ASN	CB-CG-ND2	5.70	124.95	116.40
1	A	147	SER	O-C-N	-5.32	117.05	123.27
1	A	186	GLU	CB-CG-CD	5.25	121.53	112.60
1	A	223	ASN	OD1-CG-ND2	-5.18	117.42	122.60
1	A	53[A]	VAL	CB-CA-C	-5.10	103.06	110.81
1	A	53[B]	VAL	CB-CA-C	-5.10	103.06	110.81

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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	1725	1434	1681	17	1
2	A	1	0	0	0	0
3	A	9	0	7	0	0
4	A	5	0	0	0	0
5	A	6	0	8	0	0
6	A	480	0	0	12	0
All	All	2226	1434	1696	17	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 5.

All (17) 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:145:LYS:HG3	6:A:759:HOH:O	1.22	1.34
1:A:165[B]:ASP:OD1	6:A:791:HOH:O	1.75	1.04
1:A:165[A]:ASP:HB3	6:A:787:HOH:O	1.82	0.78
1:A:145:LYS:CD	6:A:759:HOH:O	2.31	0.70
1:A:145:LYS:CG	6:A:759:HOH:O	1.99	0.70
1:A:63:ILE:HG13	6:A:723:HOH:O	1.90	0.69
1:A:145:LYS:CE	6:A:759:HOH:O	2.42	0.64
1:A:152[A]:PRO:O	6:A:447:HOH:O	2.15	0.64
1:A:60:LYS:O	6:A:723:HOH:O	2.17	0.60
1:A:145:LYS:HE3	6:A:759:HOH:O	2.03	0.57
1:A:164:SER:HB2	6:A:786:HOH:O	2.05	0.55
1:A:31[B]:VAL:CG1	1:A:65:VAL:HG13	2.37	0.55
1:A:31[B]:VAL:HG13	1:A:65:VAL:HG13	1.91	0.52
1:A:88[A]:SER:HB3	1:A:104[A]:MET:CE	2.40	0.51
1:A:87[B]:LYS:HG2	1:A:107:LYS:HB3	1.95	0.49
1:A:188(A):LYS:NZ	6:A:785:HOH:O	2.48	0.46
1:A:104[B]:MET:HE3	1:A:106:ILE:HD11	2.00	0.44

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:HH	1:A:153[A]:ASP:OD2[4_456]	1.60	0.00

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	249/223 (112%)	245 (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	A	212/184 (115%)	210 (99%)	2 (1%)	70	35

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	240[A]	GLN
1	A	240[B]	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	179	ASN

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Mol	Chain	Res	Type
1	A	210	GLN
1	A	221(A)	GLN
1	A	233	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 1 is monoatomic - leaving 3 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	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	SO4	A	253	-	4,4,4	0.79	0	6,6,6	0.94	0
3	BEN	A	252	-	9,9,9	1.89	3 (33%)	7,11,11	2.06	3 (42%)
5	GOL	A	254	-	5,5,5	0.71	0	5,5,5	0.54	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	Res	Link	Chirals	Torsions	Rings
3	BEN	A	252	-	-	4/4/4/4	0/1/1/1
5	GOL	A	254	-	-	0/4/4/4	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	252	BEN	C5-C4	-3.16	1.31	1.38
3	A	252	BEN	C6-C1	-2.85	1.35	1.39
3	A	252	BEN	C2-C1	-2.51	1.35	1.39

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	252	BEN	C5-C6-C1	3.35	123.66	120.36
3	A	252	BEN	C4-C3-C2	2.58	123.42	120.24
3	A	252	BEN	C6-C1-C2	-2.38	115.54	118.57

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	252	BEN	N2-C-C1-C2
3	A	252	BEN	N2-C-C1-C6
3	A	252	BEN	N1-C-C1-C2
3	A	252	BEN	N1-C-C1-C6

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, 95th 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>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	223/223 (100%)	-0.65	2 (0%) 81 78	3, 8, 18, 34	28 (12%)

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	148	GLY	2.7
1	A	147	SER	2.7

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 oligosaccharides 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, 95th 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	B-factors(Å ²)	Q<0.9
5	GOL	A	254	6/6	0.95	0.08	12,16,21,25	0
4	SO4	A	253	5/5	0.99	0.06	8,9,11,13	5
3	BEN	A	252	9/9	0.99	0.05	5,7,17,20	0
2	CA	A	251	1/1	1.00	0.01	5,5,5,5	0

6.5 Other polymers [i](#)

There are no such residues in this entry.