WORLDWIDE wwPDB PROTEIN DATA BANK

wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 16, 2023 – 02:42 AM EDT

PDB ID	:	1Z7X
Title	:	X-ray structure of human ribonuclease inhibitor complexed with ribonuclease I
Authors	:	McCoy, J.G.; Johnson, R.J.; Raines, R.T.; Bitto, E.; Bingman, C.A.; Wesenberg, G.E.; Allard, S.T.M.; Phillips Jr., G.N.; Center for Eukaryotic Structural Genomics (CESG)
Deposited on	:	2005-03-28
Resolution	:	1.95 Å(reported)

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

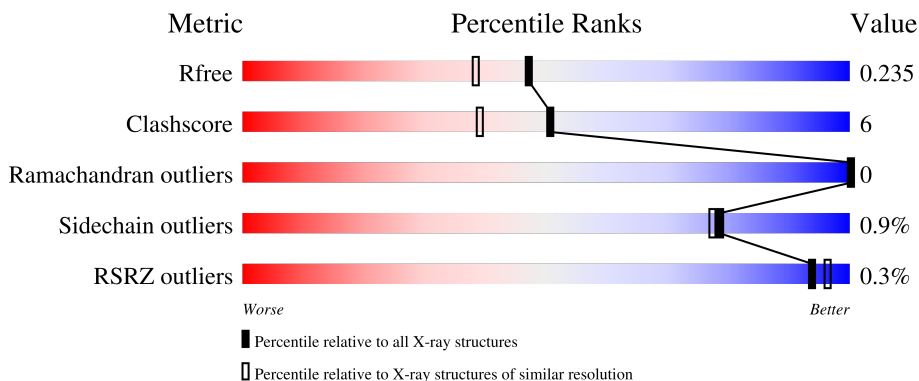|		
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:
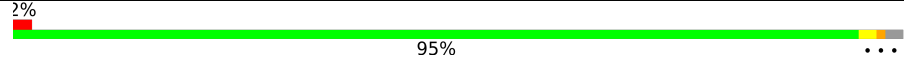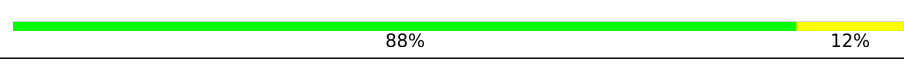*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.95 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	Percentile Ranks	Value
Rfree		0.235
Clashscore		6
Ramachandran outliers		0
Sidechain outliers		0.9%
RSRZ outliers		0.3%

Worse   *Better*

▮ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2580 (1.96-1.96)
Clashscore	141614	2705 (1.96-1.96)
Ramachandran outliers	138981	2678 (1.96-1.96)
Sidechain outliers	138945	2678 (1.96-1.96)
RSRZ outliers	127900	2539 (1.96-1.96)

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	X	129	2% ... 95% ···
1	Z	129	.% ... 83% ... 15% ·
2	W	461	88% ... 12%
2	Y	461	88% ... 11%

2 Entry composition ⓘ

There are 4 unique types of molecules in this entry. The entry contains 9795 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 Ribonuclease I.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	X	126	Total	C	N	O	S	0	0	0
			996	600	189	194	13			
1	Z	126	Total	C	N	O	S	0	0	0
			996	600	189	194	13			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
X	0	MET	-	initiating methionine	UNP P07998
Z	0	MET	-	initiating methionine	UNP P07998

- Molecule 2 is a protein called Ribonuclease inhibitor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	W	460	Total	C	N	O	S	0	0	0
			3468	2142	605	687	34			
2	Y	460	Total	C	N	O	S	0	0	0
			3468	2142	605	687	34			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
W	0	MET	-	initiating methionine	UNP P13489
Y	0	MET	-	initiating methionine	UNP P13489

- Molecule 3 is CITRIC ACID (three-letter code: CIT) (formula: $C_6H_8O_7$).

CIT

O5

O2

OH

C2 C1

HO

C6 C3

O6

C4 O1

HO

O7

C5

O O3 OH O4

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	X	1	Total 13	C 6	O 7	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	X	107	Total 107	O 107	0	0
4	W	302	Total 302	O 302	0	0
4	Z	129	Total 129	O 129	0	0
4	Y	316	Total 316	O 316	0	0

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3 Residue-property plots ⓘ

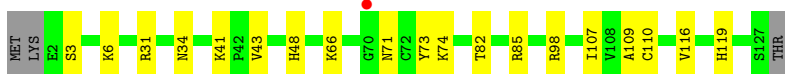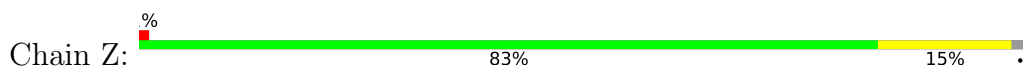These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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.
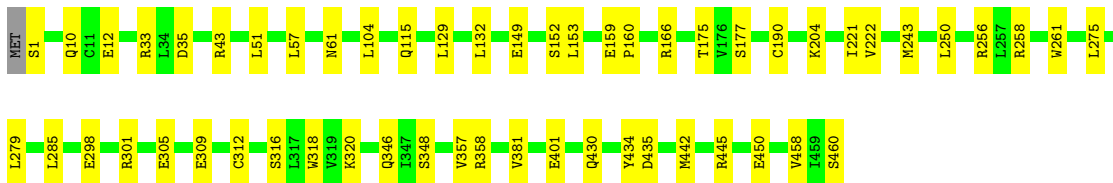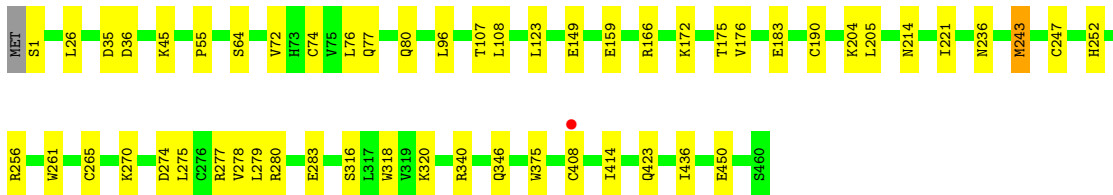
- Molecule 1: Ribonuclease I
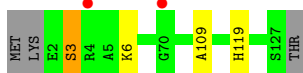
Chain X:

- Molecule 1: Ribonuclease I

Chain Z:

- Molecule 2: Ribonuclease inhibitor

Chain W:

- Molecule 2: Ribonuclease inhibitor

Chain Y:

4 Data and refinement statistics ⓘ

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	71.34Å 107.55Å 155.04Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.19 – 1.95 47.17 – 1.95	Depositor EDS
% Data completeness (in resolution range)	96.9 (47.19-1.95) 96.9 (47.17-1.95)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.82 (at 1.95Å)	Xtriage
Refinement program	REFMAC refmac_5.2.0005	Depositor
R, R_{free}	0.175 , 0.236 0.174 , 0.235	Depositor DCC
R_{free} test set	4225 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å^2)	25.9	Xtriage
Anisotropy	0.058	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 48.5	EDS
L-test for twinning[2]	$< |L| > = 0.48, < L^2 > = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	9795	wwPDB-VP
Average B, all atoms (Å^2)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.37% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.
[2]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.

5   Model quality ⓘ

5.1   Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: CIT

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	X	0.73	0/1016	0.72	0/1368
1	Z	0.78	0/1016	0.77	0/1368
2	W	0.75	0/3505	0.77	1/4744 (0.0%)
2	Y	0.74	1/3505 (0.0%)	0.77	3/4744 (0.1%)
All	All	0.75	1/9042 (0.0%)	0.76	4/12224 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	Y	247	CYS	CB-SG	-6.38	1.71	1.82

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	Y	123	LEU	CA-CB-CG	6.14	129.42	115.30
2	W	258	ARG	NE-CZ-NH1	5.77	123.18	120.30
2	Y	280	ARG	NE-CZ-NH1	5.15	122.88	120.30
2	Y	280	ARG	NE-CZ-NH2	-5.05	117.77	120.30

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	X	996	0	941	2	0
1	Z	996	0	941	21	0
2	W	3468	0	3509	45	0
2	Y	3468	0	3509	41	0
3	X	13	0	5	1	0
4	W	302	0	0	7	0
4	X	107	0	0	0	1
4	Y	316	0	0	10	0
4	Z	129	0	0	5	1
All	All	9795	0	8905	101	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 6.

The worst 5 of 101 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)
2:Y:72:VAL:HG21	2:Y:96:LEU:CD2	1.76	1.14
2:Y:72:VAL:HG21	2:Y:96:LEU:HD21	1.01	0.99
2:W:243:MET:CE	2:W:275:LEU:HD13	1.98	0.93
2:Y:72:VAL:CG2	2:Y:96:LEU:HD21	1.97	0.91
2:W:243:MET:HE2	2:W:275:LEU:HD13	1.54	0.89

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)
4:X:916:HOH:O	4:Z:257:HOH:O[2_664]	1.90	0.30

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	X	124/129 (96%)	118 (95%)	6 (5%)	0	100	100
1	Z	124/129 (96%)	121 (98%)	3 (2%)	0	100	100
2	W	458/461 (99%)	450 (98%)	8 (2%)	0	100	100
2	Y	458/461 (99%)	450 (98%)	8 (2%)	0	100	100
All	All	1164/1180 (99%)	1139 (98%)	25 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 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	X	117/120 (98%)	116 (99%)	1 (1%)	78	77
1	Z	117/120 (98%)	117 (100%)	0	100	100
2	W	396/397 (100%)	393 (99%)	3 (1%)	81	80
2	Y	396/397 (100%)	391 (99%)	5 (1%)	69	65
All	All	1026/1034 (99%)	1017 (99%)	9 (1%)	78	77

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

Mol	Chain	Res	Type
2	Y	261	TRP
2	Y	283	GLU
2	W	358	ARG
2	Y	26	LEU
2	Y	108	LEU

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

Mol	Chain	Res	Type
2	W	346	GLN
1	Z	34	ASN
1	Z	48	HIS

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Mol	Chain	Res	Type
2	Y	406	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

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	Bond lengths			Bond angles		
					Counts	RMSZ	$\#|Z| > 2$	Counts	RMSZ	$\#|Z| > 2$
3	CIT	X	900	-	12,12,12	1.36	1 (8%)	17,17,17	1.53	5 (29%)

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	CIT	X	900	-	-	2/16/16/16	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	X	900	CIT	O4-C5	-2.05	1.23	1.30

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	X	900	CIT	O6-C6-C3	2.84	117.99	113.05
3	X	900	CIT	C4-C3-C6	-2.84	104.01	110.11
3	X	900	CIT	O7-C3-C2	2.27	114.71	109.40
3	X	900	CIT	O4-C5-O3	-2.12	118.01	123.30
3	X	900	CIT	O4-C5-C4	2.02	120.85	114.35

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	X	900	CIT	O2-C1-C2-C3
3	X	900	CIT	O1-C1-C2-C3

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	X	900	CIT	1	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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(Å²)	Q<0.9
1	X	126/129 (97%)	-0.17	2 (1%)	72	79	18, 28, 42, 47	0
1	Z	126/129 (97%)	-0.35	1 (0%)	86	90	18, 24, 36, 43	0
2	W	460/461 (99%)	-0.34	0	100	100	16, 26, 38, 49	0
2	Y	460/461 (99%)	-0.22	1 (0%)	95	97	15, 27, 39, 52	0
All	All	1172/1180 (99%)	-0.27	4 (0%)	94	96	15, 26, 39, 52	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	X	70	GLY	4.3
1	Z	70	GLY	2.6
2	Y	408	CYS	2.1
1	X	4	ARG	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

6.4 Ligands ⓘ

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
3	CIT	X	900	13/13	0.95	0.13	23,29,34,36	0

6.5 Other polymers ⓘ

There are no such residues in this entry.