

wwPDB X-ray Structure Validation Summary Report (i)

Oct 23, 2021 – 08:19 AM EDT

PDB ID	:	1DBZ
Title	:	C153S MUTANT OF PEA FRUCTOSE-1,6-BISPHOSPHATASE
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Deposited on		
Resolution	:	2.65 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:

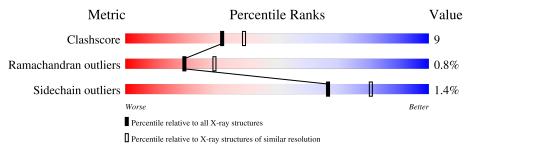
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.23.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY \, DIFFRACTION$

The reported resolution of this entry is 2.65 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	1374(2.68-2.64)
Ramachandran outliers	138981	1349 (2.68-2.64)
Sidechain outliers	138945	1349 (2.68-2.64)

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%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain		
1	А	357	68%	20%	• 11%
1	В	357	74%	16%	• 9%
1	С	357	75%	15%	• 10%
1	D	357	70%	19%	10%



 $\mathbf{2}$

Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 9686 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.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
1	Δ	317	Total	С	Ν	0	S	0	0	0
	А	317	2412	1537	398	469	8	0		0
1	В	326	Total	С	Ν	0	S	0	0	0
	D		2454	1562	403	480	9	0		
1	С	323	Total	С	Ν	0	S	0	0	0
		323	2429	1546	399	476	8	0		0
1	1 D	320	Total	С	Ν	0	S	0	0	0
			2391	1523	388	472	8		0	U

• Molecule 1 is a protein called FRUCTOSE-1,6-BISPHOSPHATASE.

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	153	SER	CYS	engineered mutation	UNP P46275
В	153	SER	CYS	engineered mutation	UNP P46275
С	153	SER	CYS	engineered mutation	UNP P46275
D	153	SER	CYS	engineered mutation	UNP P46275

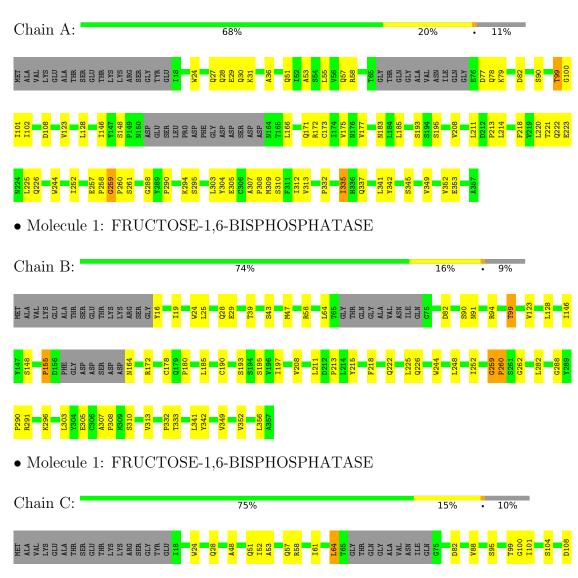




3 Residue-property plots (i)

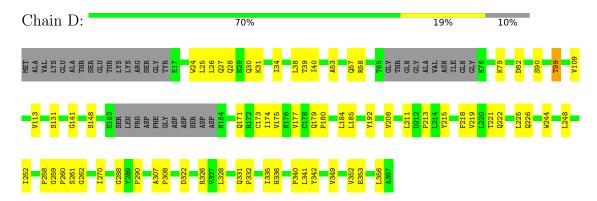
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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: FRUCTOSE-1,6-BISPHOSPHATASE

 \bullet Molecule 1: FRUCTOSE-1,6-BISPHOSPHATASE





4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source	
Space group	P 1 21 1	Depositor	
Cell constants	78.90Å 114.50Å 94.50Å	Depositor	
a, b, c, α , β , γ	90.00° 114.50° 90.00°	Depositor	
Resolution (Å)	10.00 - 2.65	Depositor	
% Data completeness	(Not available) (10.00-2.65)	Depositor	
(in resolution range)	(1000 available) (10.00-2.05)		
R_{merge}	0.10	Depositor	
R _{sym}	(Not available)	Depositor	
Refinement program	X-PLOR 3.851	Depositor	
R, R_{free}	0.235 , 0.291	Depositor	
Estimated twinning fraction	No twinning to report.	Xtriage	
Total number of atoms	9686	wwPDB-VP	
Average B, all atoms $(Å^2)$	41.0	wwPDB-VP	



5 Model quality (i)

5.1 Standard geometry (i)

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		
	Ullalli	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.48	0/2454	0.68	1/3323~(0.0%)	
1	В	0.48	0/2498	0.68	1/3389~(0.0%)	
1	С	0.47	0/2471	0.66	0/3351	
1	D	0.45	0/2433	0.66	1/3305~(0.0%)	
All	All	0.47	0/9856	0.67	3/13368~(0.0%)	

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	$\mathbf{Observed}(^{o})$	$Ideal(^{o})$
1	А	260	PRO	N-CA-CB	5.94	110.43	103.30
1	D	260	PRO	N-CA-CB	5.64	110.07	103.30
1	В	155	PRO	N-CA-CB	5.64	110.06	103.30

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.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	А	2412	0	2362	63	0
1	В	2454	0	2374	43	0
1	С	2429	0	2356	42	0
1	D	2391	0	2283	44	0
All	All	9686	0	9375	169	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 9.

The worst 5 of 169 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:175:VAL:HG22	1:A:335:ILE:HG23	1.37	1.04
1:A:214:LEU:HD22	1:C:214:LEU:HD22	1.58	0.85
1:A:304:TYR:OH	1:A:335:ILE:HB	1.78	0.84
1:A:171:GLN:HB3	1:A:335:ILE:HD11	1.61	0.83
1:A:100:GLY:HA3	1:A:166:LEU:HD21	1.65	0.78

There are no symmetry-related clashes.

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	ain Analysed Favoured Allowed		Outliers	Perce	ntiles	
1	А	311/357~(87%)	297~(96%)	12~(4%)	2(1%)	25	37
1	В	320/357~(90%)	299~(93%)	18 (6%)	3~(1%)	17	26
1	С	317/357~(89%)	295~(93%)	21 (7%)	1 (0%)	41	56
1	D	314/357~(88%)	298~(95%)	12~(4%)	4 (1%)	12	18
All	All	1262/1428~(88%)	1189 (94%)	63~(5%)	10 (1%)	19	29

5 of 10 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	А	108	ASP
1	В	260	PRO
1	D	261	SER
1	D	262	GLY
1	В	155	PRO



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 side chain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	А	259/307~(84%)	254~(98%)	5(2%)	57 74
1	В	261/307~(85%)	258~(99%)	3 (1%)	73 85
1	С	258/307~(84%)	253~(98%)	5(2%)	57 74
1	D	251/307~(82%)	250 (100%)	1 (0%)	91 95
All	All	1029/1228~(84%)	1015 (99%)	14 (1%)	67 81

5 of 14 residues with a non-rotameric sidechain are listed below:

Mol	Chain	\mathbf{Res}	Type
1	В	305	GLU
1	С	64	LEU
1	D	99	THR
1	С	296	LYS
1	С	305	GLU

Sometimes side chains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 12 such side chains are listed below:

Mol	Chain	Res	Type
1	С	240	ASN
1	С	325	GLN
1	D	325	GLN
1	D	30	GLN
1	А	325	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

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.

