



wwPDB X-ray Structure Validation Summary Report ⓘ

Jul 16, 2019 – 05:01 AM EDT

PDB ID : 2QW5
Title : CRYSTAL STRUCTURE OF A PUTATIVE SUGAR PHOSPHATE ISOMERASE/EPIMERASE (AVA4194) FROM ANABAENA VARIABILIS ATCC 29413 AT 1.78 Å RESOLUTION
Authors : Joint Center for Structural Genomics (JCSG)
Deposited on : 2007-08-09
Resolution : 1.78 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 ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.0 (224370), CSD as540be (2019)
Xtriage (Phenix) : 1.13
EDS : 2.3.2
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac : 5.8.0158
CCP4 : 7.0 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.3.2

i

X-RAY DIFFRACTION

A.

Metric	Percentile Banks	Value
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Percentile relative to all X-ray structures

Percentile relative to X-ray structures of similar resolution

Similar resolution
(#Entries, resolution range(Å))

Two horizontal bar charts showing the percentage of correct answers for two questions. The top chart shows 90% correct (green), 7% incorrect (yellow), and 3% unknown (grey). The bottom chart shows 93% correct (green), 5% incorrect (yellow), and 2% unknown (grey).

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 6066 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.

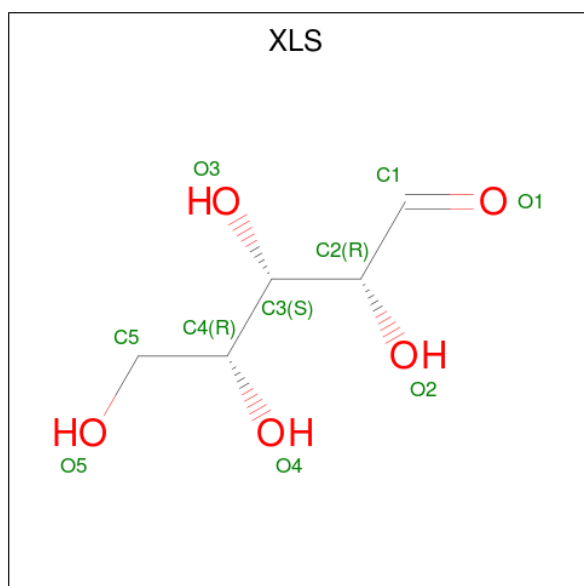
- Molecule 1 is a protein called Xylose isomerase-like TIM barrel.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	327	Total	C	N	O	Se	0	9	0
			2691	1745	444	499	3			
1	B	325	Total	C	N	O	Se	0	6	0
			2656	1723	440	490	3			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	LEADER SEQUENCE	UNP Q3M5E3
B	0	GLY	-	LEADER SEQUENCE	UNP Q3M5E3

- Molecule 2 is D-XYLOSE (LINEAR FORM) (three-letter code: XLS) (formula: C₅H₁₀O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			8	4	4		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	C	O	0	0
			8	4	4		

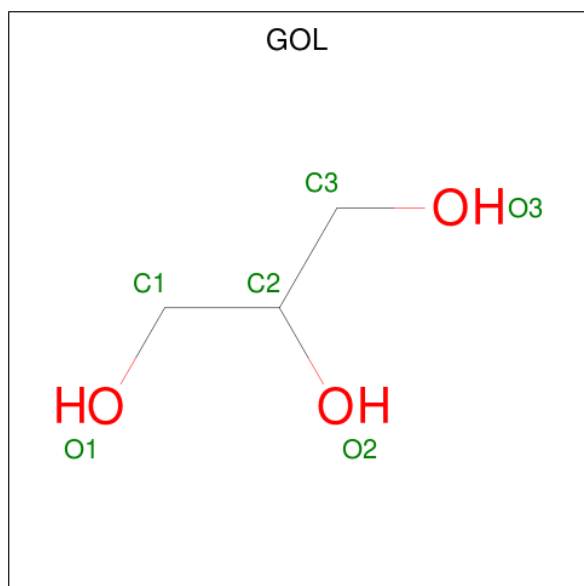
- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Zn	0	0
			1	1		
3	A	1	Total	Zn	0	0
			1	1		

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	2	Total	Cl	0	0
			2	2		
4	A	3	Total	Cl	0	0
			3	3		

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

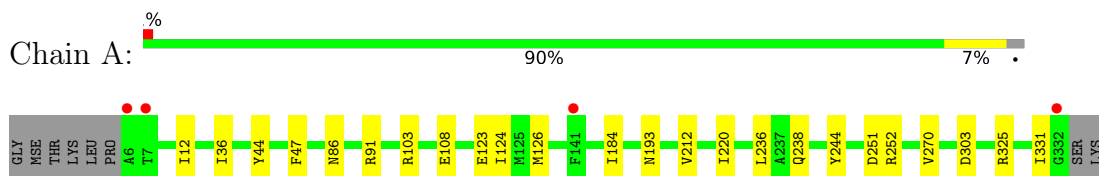
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	334	Total 334	O 334	0	0
6	B	356	Total 356	O 356	0	0

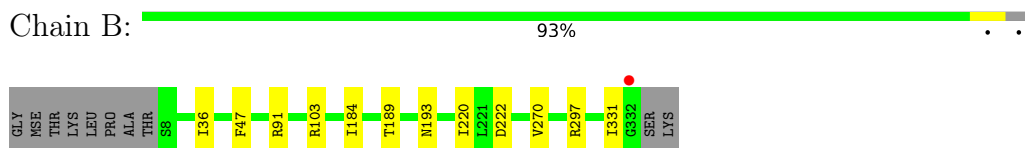
3 Residue-property plots [i](#)

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: Xylose isomerase-like TIM barrel



- Molecule 1: Xylose isomerase-like TIM barrel



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	84.24Å 84.24Å 208.17Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.11 – 1.78 42.12 – 1.78	Depositor EDS
% Data completeness (in resolution range)	99.0 (42.11-1.78) 99.0 (42.12-1.78)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.43 (at 1.78Å)	Xtriage
Refinement program	REFMAC 5.2.0019, PHENIX	Depositor
R, R_{free}	0.177 , 0.224 0.190 , 0.229	Depositor DCC
R_{free} test set	3610 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	18.2	Xtriage
Anisotropy	0.288	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 45.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6066	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.74% 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, ZN, XLS, CL

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	0.80	2/2789 (0.1%)	0.86	7/3790 (0.2%)
1	B	0.79	0/2743	0.85	3/3728 (0.1%)
All	All	0.80	2/5532 (0.0%)	0.86	10/7518 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	123	GLU	CB-CG	5.24	1.62	1.52
1	A	108	GLU	CG-CD	5.21	1.59	1.51

The worst 5 of 10 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	103	ARG	NE-CZ-NH1	8.25	124.43	120.30
1	A	252	ARG	NE-CZ-NH1	6.19	123.39	120.30
1	A	91	ARG	NE-CZ-NH1	5.99	123.29	120.30
1	A	325	ARG	NE-CZ-NH1	5.91	123.25	120.30
1	A	91	ARG	NE-CZ-NH2	-5.61	117.50	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	A	2691	0	2679	7	0
1	B	2656	0	2655	4	0
2	A	8	0	6	0	0
2	B	8	0	6	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	3	0	0	0	0
4	B	2	0	0	0	0
5	A	6	0	8	0	0
6	A	334	0	0	0	0
6	B	356	0	0	0	0
All	All	6066	0	5354	11	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 1.

The worst 5 of 11 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:B:36:ILE:HG23	1:B:47:PHE:CE2	2.40	0.56
1:A:270:VAL:HG21	1:A:331:ILE:HG21	1.91	0.53
1:A:212:VAL:HG21	1:A:236:LEU:HD13	1.92	0.51
1:B:270:VAL:HG21	1:B:331[A]:ILE:HG21	1.92	0.51
1:A:36:ILE:HG23	1:A:47:PHE:CE2	2.47	0.50

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	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	334/335 (100%)	330 (99%)	3 (1%)	1 (0%)	43	26
1	B	329/335 (98%)	325 (99%)	4 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	663/670 (99%)	655 (99%)	7 (1%)	1 (0%)	49 32

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	303	ASP

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	295/290 (102%)	292 (99%)	3 (1%)	78 71
1	B	291/290 (100%)	288 (99%)	3 (1%)	78 71
All	All	586/580 (101%)	580 (99%)	6 (1%)	76 71

5 of 6 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	251	ASP
1	B	222	ASP
1	B	189	THR
1	A	238	GLN
1	B	220	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 10 ligands modelled in this entry, 7 are 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
2	XLS	A	401	-	5,7,9	2.68	1 (20%)	6,8,11	1.31	1 (16%)
5	GOL	A	405	-	5,5,5	0.38	0	5,5,5	0.58	0
2	XLS	B	401	-	5,7,9	2.60	1 (20%)	6,8,11	1.35	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
2	XLS	A	401	-	-	4/6/8/12	-
5	GOL	A	405	-	-	2/4/4/4	-
2	XLS	B	401	-	-	4/6/8/12	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	XLS	O1-C1	5.96	1.43	1.19
2	B	401	XLS	O1-C1	5.76	1.42	1.19

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	XLS	O3-C3-C4	-2.11	104.16	109.14

There are no chirality outliers.

5 of 10 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	405	GOL	C1-C2-C3-O3
2	B	401	XLS	C1-C2-C3-C4
2	B	401	XLS	C1-C2-C3-O3
2	B	401	XLS	O2-C2-C3-C4
2	B	401	XLS	O2-C2-C3-O3

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	324/335 (96%)	-0.18	4 (1%) 79 79	14, 21, 37, 62	0
1	B	322/335 (96%)	-0.18	1 (0%) 93 93	13, 20, 36, 48	0
All	All	646/670 (96%)	-0.18	5 (0%) 86 86	13, 20, 37, 62	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	6	ALA	5.0
1	A	332	GLY	4.2
1	B	332	GLY	4.2
1	A	141[A]	PHE	2.8
1	A	7	THR	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 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, 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(\AA^2)	Q<0.9
5	GOL	A	405	6/6	0.75	0.13	46,54,55,60	0
4	CL	B	403	1/1	0.82	0.10	44,44,44,44	0
2	XLS	A	401	8/10	0.84	0.16	37,48,51,53	0
2	XLS	B	401	8/10	0.89	0.18	38,46,51,53	0
3	ZN	B	400	1/1	0.96	0.05	26,26,26,26	1
4	CL	A	404	1/1	0.96	0.11	34,34,34,34	0
4	CL	B	402	1/1	0.97	0.06	28,28,28,28	0
3	ZN	A	400	1/1	0.98	0.03	24,24,24,24	1
4	CL	A	403	1/1	0.99	0.03	20,20,20,20	0
4	CL	A	402	1/1	0.99	0.05	29,29,29,29	0

6.5 Other polymers [i](#)

There are no such residues in this entry.