



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 15, 2017 – 12:36 am GMT

PDB ID : 2QUT
Title : Dihydroxyacetone phosphate enamine intermediate in fructose-1,6-bisphosphate aldolase from rabbit muscle
Authors : St-Jean, M.; Sygusch, J.
Deposited on : 2007-08-06
Resolution : 1.88 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

<http://wwpdb.org/validation/2016/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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

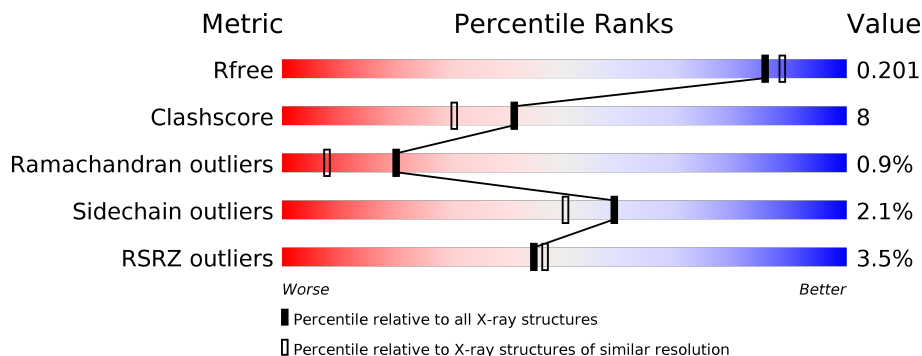
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.88 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))
R_{free}	100719	7505 (1.90-1.86)
Clashscore	112137	8369 (1.90-1.86)
Ramachandran outliers	110173	8279 (1.90-1.86)
Sidechain outliers	110143	8280 (1.90-1.86)
RSRZ outliers	101464	7571 (1.90-1.86)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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 $\leq 5\%$. 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	A	363	<div> <div>2%</div> <div> <div></div> <div>77%</div> <div>18%</div> <div>• •</div> </div> </div>
1	B	363	<div> <div>2%</div> <div> <div></div> <div>79%</div> <div>17%</div> <div>• •</div> </div> </div>
1	C	363	<div> <div>2%</div> <div> <div></div> <div>79%</div> <div>18%</div> <div>•</div> </div> </div>
1	D	363	<div> <div>7%</div> <div> <div></div> <div>78%</div> <div>21%</div> <div>•</div> </div> </div>

2 Entry composition [i](#)

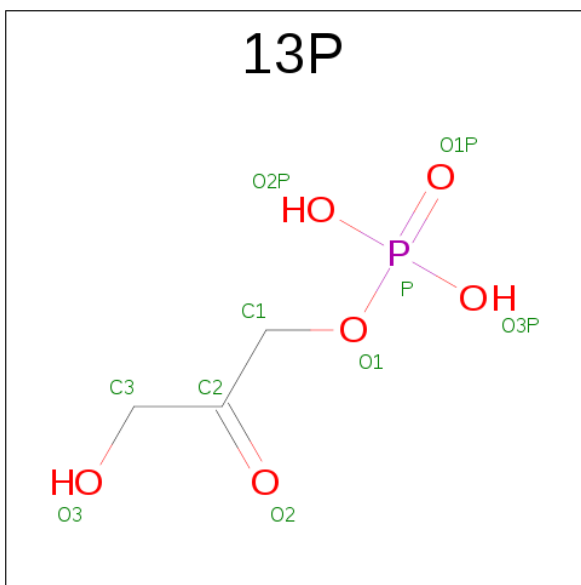
There are 3 unique types of molecules in this entry. The entry contains 13663 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Fructose-bisphosphate aldolase A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	350	Total	C	N	O	S	0	0	0
			2673	1680	475	507	11			
1	B	350	Total	C	N	O	S	0	0	0
			2673	1680	475	507	11			
1	C	353	Total	C	N	O	S	0	0	0
			2691	1690	479	511	11			
1	D	360	Total	C	N	O	S	0	0	0
			2730	1715	484	520	11			

- Molecule 2 is 1,3-DIHYDROXYACETONEPHOSPHATE (three-letter code: 13P) (formula: $C_3H_7O_6P$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	P	0	0
			9	3	5	1		
2	B	1	Total	C	O	P	0	0
			9	3	5	1		

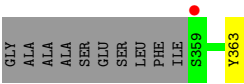
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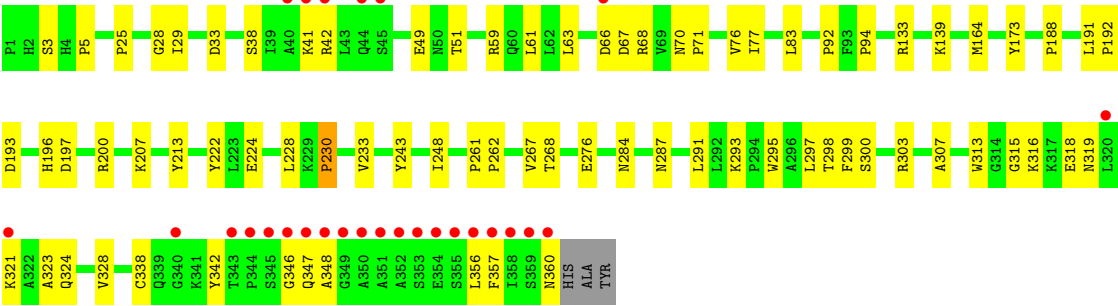
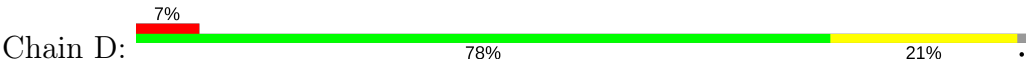
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	C	1	Total	C	O	P	0	0
			9	3	5	1		
2	D	1	Total	C	O	P	0	0
			9	3	5	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	769	Total	O	0	0
			769	769		
3	B	691	Total	O	0	0
			691	691		
3	C	715	Total	O	0	0
			715	715		
3	D	685	Total	O	0	0
			685	685		



● Molecule 1: Fructose-bisphosphate aldolase A



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	82.77Å 102.91Å 84.23Å 90.00° 98.48° 90.00°	Depositor
Resolution (Å)	50.00 – 1.88 48.17 – 1.67	Depositor EDS
% Data completeness (in resolution range)	87.1 (50.00-1.88) 78.9 (48.17-1.67)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.80 (at 1.67Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.159 , 0.191 0.168 , 0.201	Depositor DCC
R_{free} test set	5411 reflections (5.07%)	DCC
Wilson B-factor (Å ²)	17.7	Xtriage
Anisotropy	0.558	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 84.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.017 for l,-k,h	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	13663	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 5.36% 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: 13P

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	0.29	0/2725	0.58	1/3691 (0.0%)
1	B	0.29	0/2725	0.58	0/3691
1	C	0.30	0/2743	0.58	0/3715
1	D	0.29	0/2782	0.57	0/3770
All	All	0.29	0/10975	0.58	1/14867 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1
1	C	0	1
1	D	0	1
All	All	0	3

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	300	SER	N-CA-C	-5.29	96.72	111.00

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	213	TYR	Sidechain
1	C	213	TYR	Sidechain

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Mol	Chain	Res	Type	Group
1	D	213	TYR	Sidechain

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	2673	0	2693	46	0
1	B	2673	0	2692	44	0
1	C	2691	0	2708	44	0
1	D	2730	0	2753	52	0
2	A	9	0	5	1	0
2	B	9	0	5	0	0
2	C	9	0	5	0	0
2	D	9	0	5	0	0
3	A	769	0	0	9	0
3	B	691	0	0	6	0
3	C	715	0	0	7	0
3	D	685	0	0	10	0
All	All	13663	0	10866	175	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 8.

The worst 5 of 175 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:193:ASP:HA	1:B:237:HIS:HD2	1.42	0.83
1:B:193:ASP:HA	1:B:237:HIS:CD2	2.24	0.72
1:B:291:LEU:O	1:B:293:LYS:HD3	1.90	0.72
1:A:155:GLU:HG2	3:A:3624:HOH:O	1.91	0.71
1:C:336:LEU:HD11	1:C:347:GLN:HA	1.72	0.70

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	346/363 (95%)	329 (95%)	14 (4%)	3 (1%)	20	8
1	B	346/363 (95%)	329 (95%)	14 (4%)	3 (1%)	20	8
1	C	349/363 (96%)	335 (96%)	12 (3%)	2 (1%)	28	15
1	D	358/363 (99%)	343 (96%)	11 (3%)	4 (1%)	17	5
All	All	1399/1452 (96%)	1336 (96%)	51 (4%)	12 (1%)	20	8

5 of 12 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	5	PRO
1	A	188	PRO
1	B	5	PRO
1	B	188	PRO
1	C	5	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 sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	284/291 (98%)	276 (97%)	8 (3%)	49	37
1	B	284/291 (98%)	278 (98%)	6 (2%)	59	50
1	C	285/291 (98%)	280 (98%)	5 (2%)	64	57
1	D	289/291 (99%)	284 (98%)	5 (2%)	66	59
All	All	1142/1164 (98%)	1118 (98%)	24 (2%)	59	50

5 of 24 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	193	ASP
1	B	295	TRP
1	D	230	PRO
1	B	230	PRO
1	B	277	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 28 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	220	HIS
1	C	2	HIS
1	D	284	ASN
1	B	237	HIS
1	B	319	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

4 ligands are 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
2	13P	A	3001	1	8,8,9	1.11	1 (12%)	10,10,12	1.07	1 (10%)
2	13P	B	3002	1	8,8,9	1.08	1 (12%)	10,10,12	1.12	1 (10%)
2	13P	C	3003	1	8,8,9	1.15	1 (12%)	10,10,12	1.12	1 (10%)
2	13P	D	3004	1	8,8,9	1.14	1 (12%)	10,10,12	1.12	1 (10%)

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	13P	A	3001	1	-	0/6/6/8	0/0/0/0
2	13P	B	3002	1	-	0/6/6/8	0/0/0/0
2	13P	C	3003	1	-	0/6/6/8	0/0/0/0
2	13P	D	3004	1	-	0/6/6/8	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	3003	13P	C2-C3	-2.69	1.35	1.50
2	D	3004	13P	C2-C3	-2.66	1.35	1.50
2	A	3001	13P	C2-C3	-2.65	1.35	1.50
2	B	3002	13P	C2-C3	-2.58	1.35	1.50

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	3001	13P	C1-C2-C3	2.91	120.61	113.86
2	B	3002	13P	C1-C2-C3	3.11	121.06	113.86
2	D	3004	13P	C1-C2-C3	3.13	121.10	113.86
2	C	3003	13P	C1-C2-C3	3.14	121.12	113.86

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	3001	13P	1	0

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	350/363 (96%)	-0.18	7 (2%) 65 68	10, 19, 43, 64	2 (0%)
1	B	350/363 (96%)	-0.16	6 (1%) 70 73	9, 19, 45, 67	2 (0%)
1	C	353/363 (97%)	-0.08	9 (2%) 58 60	9, 18, 41, 71	4 (1%)
1	D	360/363 (99%)	0.18	27 (7%) 15 16	11, 23, 60, 106	0
All	All	1413/1452 (97%)	-0.06	49 (3%) 44 46	9, 20, 47, 106	8 (0%)

The worst 5 of 49 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	345	SER	18.7
1	A	345	SER	10.2
1	C	348	ALA	9.9
1	A	359	SER	9.4
1	D	349	GLY	8.4

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
2	13P	C	3003	9/10	0.96	0.12	0.86	21,22,23,23	0
2	13P	B	3002	9/10	0.96	0.11	0.60	23,23,25,29	0
2	13P	D	3004	9/10	0.95	0.11	0.43	26,28,29,30	0
2	13P	A	3001	9/10	0.97	0.08	0.41	19,21,22,23	0

6.5 Other polymers [i](#)

There are no such residues in this entry.