



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 11:51 pm BST

PDB ID : 1TKS
Title : Crystal structure of 3,4-Dihydroxy-2-butanone 4-phosphate Synthase of *Candida albicans*
Authors : Echt, S.; Bauer, S.; Steinbacher, S.; Huber, R.; Bacher, A.; Fischer, M.
Deposited on : 2004-06-09
Resolution : 1.60 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
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)	:	2.11

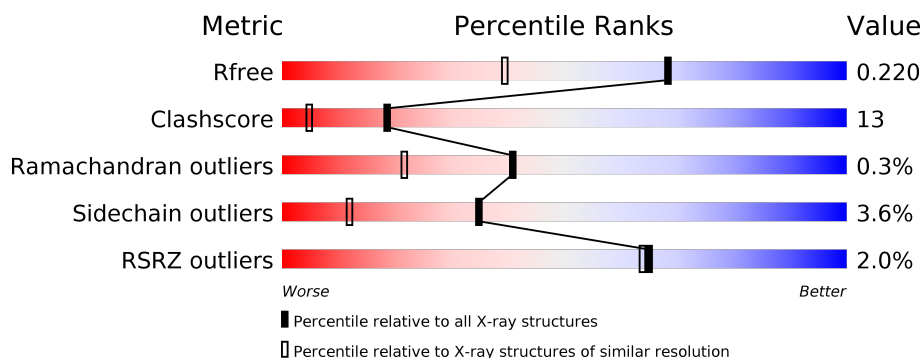
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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}	130704	3398 (1.60-1.60)
Clashscore	141614	3665 (1.60-1.60)
Ramachandran outliers	138981	3564 (1.60-1.60)
Sidechain outliers	138945	3563 (1.60-1.60)
RSRZ outliers	127900	3321 (1.60-1.60)

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 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 $\leq 5\%$. 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	204	<div> <div>%</div> <div> <div></div> <div>77%</div> <div>17%</div> <div>• •</div> </div> </div>
1	B	204	<div> <div>3%</div> <div> <div></div> <div>75%</div> <div>19%</div> <div>• •</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3710 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.

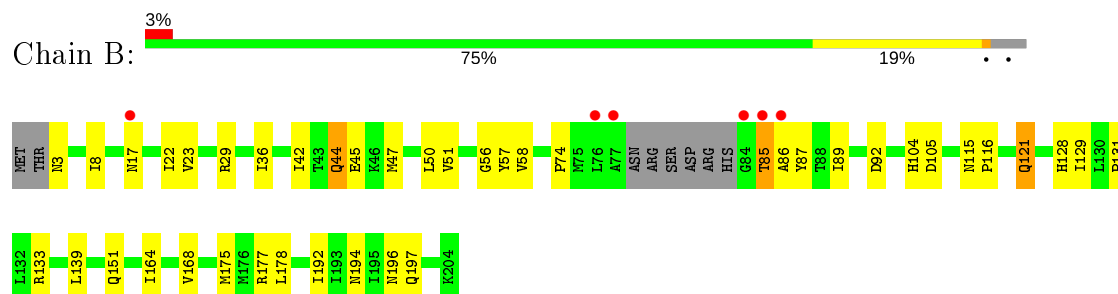
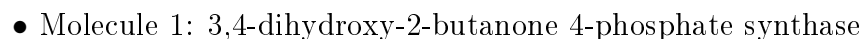
- Molecule 1 is a protein called 3,4-dihydroxy-2-butanone 4-phosphate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	196	Total	C	N	O	S	11	0	0
			1517	955	259	293	10			
1	B	196	Total	C	N	O	S	6	0	0
			1517	955	259	293	10			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	328	Total	O	0	0
			328	328		
2	B	348	Total	O	0	0
			348	348		

- Molecule 1: 3,4-dihydroxy-2-butanone 4-phosphate synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	40.44Å 48.13Å 59.80Å 66.23° 71.98° 89.86°	Depositor
Resolution (Å)	20.00 – 1.60 19.56 – 1.60	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-1.60) 84.4 (19.56-1.60)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.60 (at 1.60Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.192 , 0.219 0.193 , 0.220	Depositor DCC
R_{free} test set	4367 reflections (10.09%)	wwPDB-VP
Wilson B-factor (Å ²)	13.8	Xtriage
Anisotropy	0.277	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 44.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.018 for h,-k,h-l 0.476 for -h,k,k-l 0.017 for -h,-k,-h-k+l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3710	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.16% 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

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	0/1540	0.58	0/2080
1	B	0.29	0/1540	0.58	0/2080
All	All	0.29	0/3080	0.58	0/4160

There are no bond length outliers.

There are no bond angle outliers.

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	1517	0	1529	46	0
1	B	1517	0	1529	36	0
2	A	328	0	0	7	0
2	B	348	0	0	3	0
All	All	3710	0	3058	80	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 13.

All (80) 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:47:MET:HE2	1:B:51:VAL:HG23	1.32	1.10

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:47:MET:HE3	1:B:50:LEU:HB3	1.62	0.80
1:B:29:ARG:HD3	1:B:178:LEU:HD23	1.62	0.80
1:A:115:ASN:HD22	1:A:117:ASN:H	1.33	0.77
1:A:8:ILE:HD13	1:A:201:TYR:HB2	1.70	0.72
1:B:121:GLN:NE2	1:B:121:GLN:H	1.88	0.71
1:A:36:ILE:HG12	1:A:164:ILE:HG22	1.73	0.71
2:A:515:HOH:O	1:B:89:ILE:HG12	1.94	0.67
1:A:22:ILE:HD12	1:A:192:ILE:CD1	2.25	0.66
1:A:8:ILE:HD11	1:A:197:GLN:HG2	1.77	0.66
1:B:121:GLN:HE21	1:B:121:GLN:H	1.42	0.66
1:B:8:ILE:HD11	1:B:197:GLN:HG2	1.78	0.65
1:A:22:ILE:HD12	1:A:192:ILE:HD12	1.79	0.64
1:B:85:THR:HA	1:B:87:TYR:CE1	2.36	0.60
1:A:194:ASN:ND2	1:A:197:GLN:H	2.00	0.60
1:A:115:ASN:ND2	1:A:117:ASN:H	1.99	0.60
1:A:106:ARG:HA	1:A:129:ILE:HD11	1.83	0.60
1:A:121:GLN:NE2	1:A:121:GLN:H	1.98	0.60
1:A:106:ARG:HG2	1:A:129:ILE:HD11	1.84	0.60
1:A:139:LEU:H	1:A:151:GLN:NE2	2.01	0.58
1:A:56:GLY:HA3	1:B:56:GLY:HA3	1.85	0.58
1:B:36:ILE:HG23	1:B:164:ILE:HG22	1.85	0.58
1:B:104:HIS:HD2	1:B:105:ASP:OD1	1.86	0.58
1:B:47:MET:HE2	1:B:51:VAL:CG2	2.22	0.58
1:A:22:ILE:HB	1:A:192:ILE:HD12	1.86	0.57
1:B:47:MET:HE1	1:B:58:VAL:HG21	1.87	0.55
1:B:36:ILE:HG12	1:B:164:ILE:HG22	1.88	0.54
1:A:8:ILE:HD13	1:A:201:TYR:CB	2.37	0.53
1:B:139:LEU:HB2	1:B:151:GLN:HE21	1.74	0.51
1:B:23:VAL:HG21	1:B:36:ILE:HD12	1.93	0.51
1:B:177:ARG:HB2	2:B:242:HOH:O	2.10	0.50
1:B:139:LEU:H	1:B:151:GLN:NE2	2.08	0.50
1:B:8:ILE:CD1	1:B:197:GLN:HG2	2.40	0.50
1:B:121:GLN:N	1:B:121:GLN:HE21	2.10	0.50
1:A:191:LYS:C	1:A:192:ILE:HD13	2.33	0.49
1:A:23:VAL:HG21	1:A:36:ILE:HD12	1.94	0.49
1:A:42:ILE:HD12	1:A:47:MET:HB2	1.95	0.48
1:A:24:MET:CE	1:A:178:LEU:HA	2.44	0.48
1:A:42:ILE:HD11	1:A:47:MET:N	2.29	0.48
1:A:36:ILE:HG12	1:A:164:ILE:CG2	2.43	0.48
1:A:8:ILE:HD12	2:A:263:HOH:O	2.13	0.48
1:A:188:HIS:HE1	2:A:429:HOH:O	1.97	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:8:ILE:CD1	1:A:201:TYR:HB2	2.41	0.47
1:A:141:LYS:HD2	2:A:461:HOH:O	2.14	0.47
1:B:74:PRO:HD3	1:B:133:ARG:HH11	1.79	0.47
1:A:4:ILE:HG23	1:A:182:ILE:CD1	2.45	0.47
1:A:128:HIS:C	1:A:129:ILE:HD12	2.35	0.47
1:B:92:ASP:OD1	1:B:128:HIS:HD2	1.97	0.46
1:A:4:ILE:HG23	1:A:182:ILE:HD11	1.98	0.46
1:B:129:ILE:O	1:B:131:PRO:HD3	2.16	0.46
1:A:129:ILE:O	1:A:131:PRO:HD3	2.16	0.46
1:A:24:MET:HE2	1:A:178:LEU:HD12	1.97	0.46
1:B:3:ASN:HB2	2:B:356:HOH:O	2.15	0.46
1:B:85:THR:HA	1:B:87:TYR:CD1	2.51	0.46
1:A:139:LEU:HB2	1:A:151:GLN:NE2	2.31	0.45
1:A:139:LEU:HB2	1:A:151:GLN:HE21	1.81	0.45
1:A:131:PRO:O	1:A:132:LEU:HD23	2.17	0.44
1:B:139:LEU:HB2	1:B:151:GLN:NE2	2.32	0.44
1:B:168:VAL:HG23	2:B:258:HOH:O	2.16	0.44
1:B:194:ASN:H	1:B:197:GLN:HE21	1.66	0.44
1:A:8:ILE:HD11	1:A:197:GLN:O	2.17	0.44
1:A:115:ASN:C	1:A:115:ASN:HD22	2.21	0.43
1:A:182:ILE:HD12	2:A:265:HOH:O	2.18	0.43
1:B:8:ILE:HD11	1:B:197:GLN:CG	2.48	0.43
1:A:64:GLU:HG3	1:A:134:ALA:O	2.18	0.43
1:B:42:ILE:HG23	1:B:42:ILE:O	2.19	0.42
1:A:84:GLY:HA2	2:A:425:HOH:O	2.19	0.42
1:A:106:ARG:HG2	1:A:129:ILE:CD1	2.49	0.42
1:A:37:MET:SD	1:A:42:ILE:HG12	2.60	0.42
1:B:44:GLN:HE21	1:B:45:GLU:N	2.17	0.41
1:A:188:HIS:HD2	2:A:497:HOH:O	2.04	0.41
1:A:24:MET:HE2	1:A:178:LEU:HA	2.03	0.41
1:B:115:ASN:HA	1:B:116:PRO:HD3	1.93	0.41
1:A:29:ARG:HG2	1:A:29:ARG:NH2	2.36	0.41
1:B:22:ILE:HB	1:B:192:ILE:HG12	2.02	0.40
1:B:85:THR:HG23	1:B:86:ALA:N	2.36	0.40
1:B:8:ILE:N	1:B:8:ILE:HD12	2.36	0.40
1:A:42:ILE:O	1:A:42:ILE:CG2	2.70	0.40
1:A:101:ILE:HD11	1:B:57:TYR:CB	2.52	0.40
1:A:29:ARG:HH21	1:A:29:ARG:HG2	1.86	0.40

There are no symmetry-related clashes.

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	A	192/204 (94%)	184 (96%)	8 (4%)	0	100	100
1	B	192/204 (94%)	184 (96%)	7 (4%)	1 (0%)	29	11
All	All	384/408 (94%)	368 (96%)	15 (4%)	1 (0%)	41	21

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	85	THR

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	A	165/173 (95%)	158 (96%)	7 (4%)	30	9
1	B	165/173 (95%)	160 (97%)	5 (3%)	41	16
All	All	330/346 (95%)	318 (96%)	12 (4%)	35	12

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	42	ILE
1	A	44	GLN
1	A	115	ASN
1	A	121	GLN
1	A	151	GLN

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Mol	Chain	Res	Type
1	A	175	MET
1	A	177	ARG
1	B	17	ASN
1	B	44	GLN
1	B	121	GLN
1	B	175	MET
1	B	196	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	ASN
1	A	44	GLN
1	A	115	ASN
1	A	117	ASN
1	A	121	GLN
1	A	151	GLN
1	A	183	GLN
1	A	188	HIS
1	A	194	ASN
1	B	3	ASN
1	B	31	ASN
1	B	44	GLN
1	B	104	HIS
1	B	117	ASN
1	B	121	GLN
1	B	128	HIS
1	B	151	GLN
1	B	197	GLN

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 [i](#)

There are no carbohydrates 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](#)

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	195/204 (95%)	-0.28	2 (1%) 82 82	8, 11, 21, 29	1 (0%)
1	B	196/204 (96%)	-0.23	6 (3%) 49 46	7, 11, 21, 37	1 (0%)
All	All	391/408 (95%)	-0.26	8 (2%) 65 64	7, 11, 21, 37	2 (0%)

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	85	THR	5.9
1	B	77	ALA	3.0
1	A	202	ILE	2.9
1	B	86	ALA	2.8
1	A	17	ASN	2.4
1	B	17	ASN	2.1
1	B	84	GLY	2.1
1	B	76	LEU	2.0

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.