Full wwPDB X-ray Structure Validation Report ⓘ

Jun 12, 2024 – 09:21 PM EDT

PDB ID	:	3HIN
Title	:	CRYSTAL STRUCTURE OF putative enoyl-CoA hydratase from Rhodopseudomonas palustris CGA009
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Deposited on	:	2009-05-20
Resolution	:	2.00 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 types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

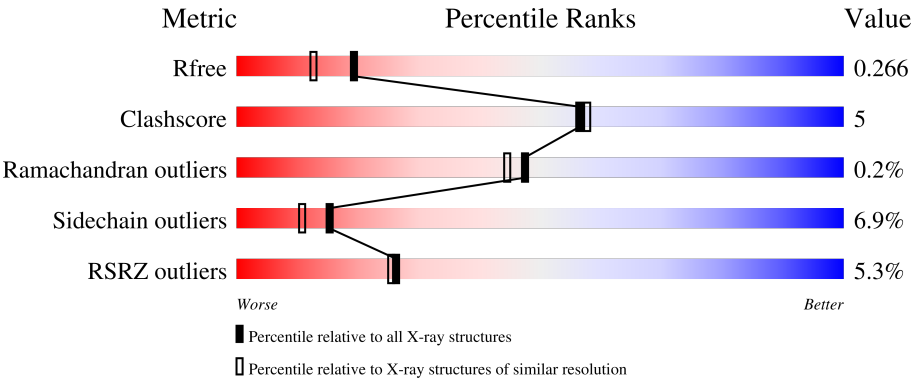
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.20.1
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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	A	275	5% 81% 11% • 6%
1	B	275	5% 77% 11% • 9%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 4045 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 Putative 3-hydroxybutyryl-CoA dehydratase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	258	Total	C	N	O	S	0	0	0
			1925	1210	347	358	10			
1	B	251	Total	C	N	O	S	0	0	0
			1878	1180	338	350	10			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP Q6N8W7
A	2	SER	-	expression tag	UNP Q6N8W7
A	268	GLU	-	expression tag	UNP Q6N8W7
A	269	GLY	-	expression tag	UNP Q6N8W7
A	270	HIS	-	expression tag	UNP Q6N8W7
A	271	HIS	-	expression tag	UNP Q6N8W7
A	272	HIS	-	expression tag	UNP Q6N8W7
A	273	HIS	-	expression tag	UNP Q6N8W7
A	274	HIS	-	expression tag	UNP Q6N8W7
A	275	HIS	-	expression tag	UNP Q6N8W7
B	1	MET	-	expression tag	UNP Q6N8W7
B	2	SER	-	expression tag	UNP Q6N8W7
B	268	GLU	-	expression tag	UNP Q6N8W7
B	269	GLY	-	expression tag	UNP Q6N8W7
B	270	HIS	-	expression tag	UNP Q6N8W7
B	271	HIS	-	expression tag	UNP Q6N8W7
B	272	HIS	-	expression tag	UNP Q6N8W7
B	273	HIS	-	expression tag	UNP Q6N8W7
B	274	HIS	-	expression tag	UNP Q6N8W7
B	275	HIS	-	expression tag	UNP Q6N8W7

- Molecule 2 is water.

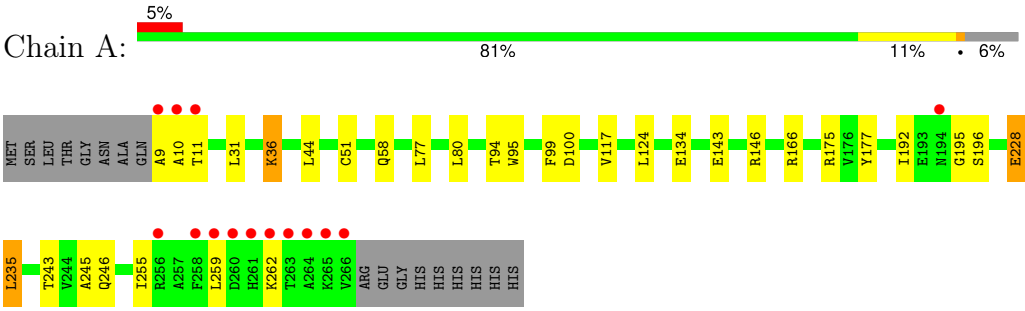
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	131	Total O 131 131	0	0
2	B	111	Total O 111 111	0	0

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 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.

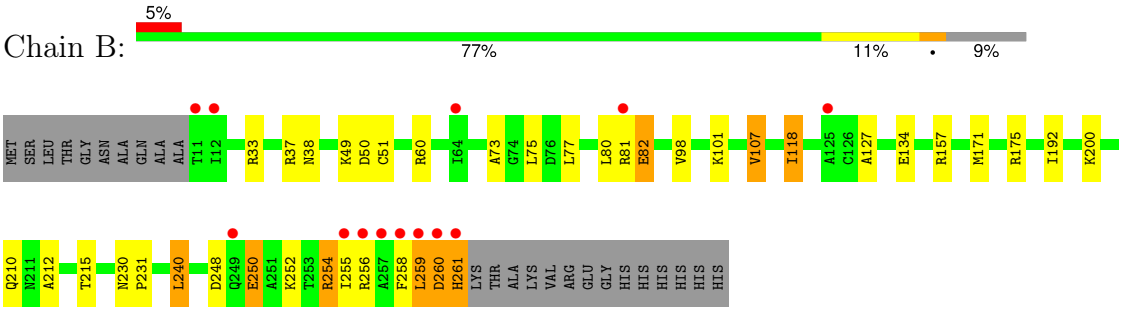
• Molecule 1: Putative 3-hydroxybutyryl-CoA dehydratase

Chain A:



• Molecule 1: Putative 3-hydroxybutyryl-CoA dehydratase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants a, b, c, α, β, γ	124.11Å 124.11Å 124.11Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.25 – 2.00 29.25 – 2.00	Depositor EDS
% Data completeness (in resolution range)	100.0 (29.25-2.00) 88.6 (29.25-2.00)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.56 (at 2.01Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.225 , 0.274 0.221 , 0.266	Depositor DCC
R_{free} test set	1929 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å^2)	35.4	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 40.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.35$	Xtriage
Estimated twinning fraction	0.039 for l,-k,h	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	4045	wwPDB-VP
Average B, all atoms (Å^2)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.44% of the height of the origin peak. No significant pseudotranslation is detected.*

[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

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	1.06	5/1955 (0.3%)	0.94	4/2651 (0.2%)
1	B	1.04	2/1908 (0.1%)	0.98	3/2588 (0.1%)
All	All	1.05	7/3863 (0.2%)	0.96	7/5239 (0.1%)

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	A	0	1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	134	GLU	CB-CG	-8.51	1.35	1.52
1	A	117	VAL	CB-CG2	6.96	1.67	1.52
1	A	51	CYS	CB-SG	-6.32	1.71	1.82
1	A	228	GLU	CG-CD	6.30	1.61	1.51
1	B	51	CYS	CB-SG	-5.32	1.73	1.81
1	A	245	ALA	CA-CB	-5.17	1.41	1.52
1	A	134	GLU	CB-CG	-5.02	1.42	1.52

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	240	LEU	CB-CG-CD1	9.85	127.74	111.00
1	A	175	ARG	NE-CZ-NH1	-6.27	117.17	120.30
1	A	100	ASP	CB-CG-OD2	5.98	123.68	118.30
1	B	171	MET	CG-SD-CE	-5.52	91.36	100.20
1	B	37	ARG	NE-CZ-NH2	5.51	123.05	120.30
1	A	235	LEU	CA-CB-CG	5.49	127.92	115.30
1	A	235	LEU	CB-CG-CD1	5.05	119.59	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	195	GLY	Peptide

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	A	1925	0	1961	13	0
1	B	1878	0	1904	26	0
2	A	131	0	0	3	0
2	B	111	0	0	0	0
All	All	4045	0	3865	38	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 5.

All (38) 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:252:LYS:O	1:B:255:ILE:HG22	1.76	0.85
1:B:260:ASP:O	1:B:261:HIS:HB2	1.80	0.81
1:A:228:GLU:OE2	2:A:284:HOH:O	2.13	0.66
1:B:230:ASN:HD22	1:B:231:PRO:HD2	1.63	0.63
1:B:210:GLN:O	1:B:254:ARG:NH1	2.27	0.63
1:B:77:LEU:O	1:B:80:LEU:HB2	2.00	0.61
1:B:77:LEU:HA	1:B:80:LEU:HD13	1.84	0.60
1:B:192:ILE:HD11	1:B:200:LYS:HG2	1.82	0.59
1:A:99:PHE:CE1	1:A:124:LEU:HD13	2.38	0.59
1:B:230:ASN:ND2	1:B:231:PRO:HD2	2.19	0.57
1:B:212:ALA:O	1:B:215:THR:HB	2.06	0.56
1:B:248:ASP:O	1:B:252:LYS:HG2	2.08	0.54
1:B:260:ASP:O	1:B:261:HIS:CB	2.56	0.54
1:A:99:PHE:HE1	1:A:124:LEU:HD13	1.74	0.53
1:A:9:ALA:HB1	2:A:373:HOH:O	2.10	0.52
1:A:94:THR:HG22	1:A:95:TRP:HD1	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:230:ASN:HD22	1:B:231:PRO:CD	2.24	0.50
1:B:60:ARG:HA	1:B:107:VAL:HG21	1.94	0.49
1:A:243:THR:HA	1:A:246:GLN:HG3	1.95	0.48
1:A:77:LEU:O	1:A:80:LEU:CD1	2.60	0.48
1:B:49:LYS:HB2	1:B:98:VAL:HG22	1.96	0.47
1:B:82:GLU:OE1	1:B:82:GLU:HA	2.14	0.47
1:A:44:LEU:C	1:A:44:LEU:HD23	2.36	0.46
1:B:73:ALA:O	1:B:118:ILE:HG22	2.15	0.46
1:A:9:ALA:CB	2:A:373:HOH:O	2.64	0.45
1:B:250:GLU:OE1	1:B:254:ARG:HD2	2.16	0.45
1:B:75:LEU:HA	1:B:75:LEU:HD12	1.66	0.45
1:B:60:ARG:O	1:B:107:VAL:HG22	2.17	0.44
1:B:252:LYS:C	1:B:255:ILE:HG22	2.36	0.44
1:A:166:ARG:NH2	1:A:177:TYR:OH	2.48	0.44
1:A:230:ASN:OD1	1:B:230:ASN:ND2	2.53	0.42
1:A:230:ASN:ND2	1:A:233:THR:OG1	2.53	0.42
1:B:33:ARG:HB2	1:B:38:ASN:HA	2.02	0.42
1:B:127:ALA:HA	1:B:157:ARG:HD3	2.03	0.41
1:A:36:LYS:H	1:A:36:LYS:HG2	1.75	0.41
1:B:255:ILE:O	1:B:259:LEU:HB2	2.20	0.41
1:B:252:LYS:HA	1:B:255:ILE:HG22	2.04	0.40
1:B:175:ARG:HH11	1:B:175:ARG:HD3	1.72	0.40

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	256/275 (93%)	249 (97%)	6 (2%)	1 (0%)	34 30
1	B	249/275 (90%)	240 (96%)	9 (4%)	0	100 100
All	All	505/550 (92%)	489 (97%)	15 (3%)	1 (0%)	47 44

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	10	ALA

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	198/212 (93%)	185 (93%)	13 (7%)	16 12
1	B	194/212 (92%)	180 (93%)	14 (7%)	14 9
All	All	392/424 (92%)	365 (93%)	27 (7%)	15 11

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	THR
1	A	31	LEU
1	A	36	LYS
1	A	58	GLN
1	A	143	GLU
1	A	146	ARG
1	A	192	ILE
1	A	196	SER
1	A	230	ASN
1	A	235	LEU
1	A	255	ILE
1	A	259	LEU
1	A	262	LYS
1	B	50	ASP
1	B	81	ARG
1	B	82	GLU
1	B	101	LYS
1	B	107	VAL
1	B	118	ILE
1	B	240	LEU
1	B	250	GLU
1	B	254	ARG

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Mol	Chain	Res	Type
1	B	256	ARG
1	B	258	PHE
1	B	259	LEU
1	B	260	ASP
1	B	261	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	58	GLN
1	A	230	ASN
1	A	232	GLN
1	B	230	ASN
1	B	232	GLN
1	B	249	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

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, 95^{th} 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	258/275 (93%)	-0.13	14 (5%)	25	24	25, 36, 63, 104	0
1	B	251/275 (91%)	-0.07	13 (5%)	27	26	26, 36, 64, 100	0
All	All	509/550 (92%)	-0.10	27 (5%)	26	25	25, 36, 64, 104	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	266	VAL	9.7
1	A	264	ALA	9.5
1	A	263	THR	7.8
1	B	260	ASP	6.4
1	B	261	HIS	5.7
1	B	258	PHE	5.6
1	B	259	LEU	5.4
1	A	265	LYS	5.3
1	A	9	ALA	5.2
1	A	261	HIS	4.5
1	A	10	ALA	4.2
1	B	11	THR	3.6
1	A	262	LYS	3.4
1	B	249	GLN	3.2
1	B	257	ALA	3.0
1	A	260	ASP	2.9
1	B	12	ILE	2.7
1	A	258	PHE	2.5
1	B	125	ALA	2.4
1	B	81	ARG	2.4
1	B	256	ARG	2.3
1	A	11	THR	2.3
1	A	256	ARG	2.3
1	A	194	ASN	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	255	ILE	2.2
1	A	259	LEU	2.1
1	B	64	ILE	2.1

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 monosaccharides 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.