



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 17, 2023 – 09:10 AM EST

PDB ID : 2GLZ
Title : Crystal structure of a formylmethanofuran dehydrogenase subunit e-like protein (dhaf_2992) from desulfitobacterium hafniense dcb-2 at 1.45 Å resolution
Authors : Joint Center for Structural Genomics (JCSG)
Deposited on : 2006-04-05
Resolution : 1.45 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
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.31.2
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) : 2.31.2

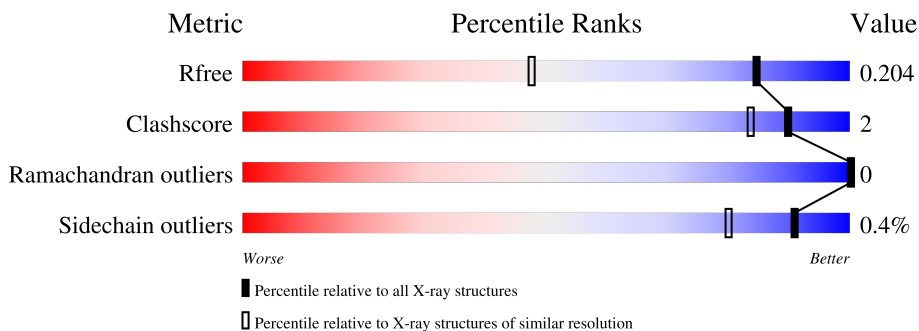
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.45 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	1156 (1.46-1.46)
Clashscore	141614	1202 (1.46-1.46)
Ramachandran outliers	138981	1178 (1.46-1.46)
Sidechain outliers	138945	1178 (1.46-1.46)

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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	153	93% (with small segments for 1, 2, 3 types of outliers and unmodelled residues)
1	B	153	93% (with small segments for 1, 2, 3 types of outliers and unmodelled residues)

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 2949 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 similar to Formylmethanofuran dehydrogenase subunit E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	N	O	S				Se
1	A	149	1215	780	211	217	4	3	0	8	0
1	B	148	1227	783	215	222	4	3	0	9	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	GB 68208641
A	1	MSE	MET	modified residue	GB 68208641
A	36	MSE	MET	modified residue	GB 68208641
A	82	MSE	MET	modified residue	GB 68208641
B	0	GLY	-	expression tag	GB 68208641
B	1	MSE	MET	modified residue	GB 68208641
B	36	MSE	MET	modified residue	GB 68208641
B	82	MSE	MET	modified residue	GB 68208641

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 3 is NICKEL (II) ION (three-letter code: NI) (formula: Ni).

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

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



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

Continued on next page...

Continued from previous page...

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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	214	Total O 216 216	0	2
5	B	214	Total O 215 215	0	1

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.

- Molecule 1: similar to Formylmethanofuran dehydrogenase subunit E

Chain A:  93%



- Molecule 1: similar to Formylmethanofuran dehydrogenase subunit E

Chain B:  93%



4 Data and refinement statistics i

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	46.42Å 84.79Å 100.71Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	27.20 – 1.45 27.20 – 1.45	Depositor EDS
% Data completeness (in resolution range)	99.9 (27.20-1.45) 99.9 (27.20-1.45)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.81 (at 1.45Å)	Xtrriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.171 , 0.198 0.177 , 0.204	Depositor DCC
R_{free} test set	3593 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	17.1	Xtrriage
Anisotropy	0.636	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 42.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2949	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 37.71 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.1174e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: EDO, NI, ZN

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	0.71	0/1262	0.81	0/1710
1	B	0.75	0/1272	0.80	1/1721 (0.1%)
All	All	0.73	0/2534	0.81	1/3431 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	27	ARG	NE-CZ-NH1	-5.40	117.60	120.30

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	A	1215	0	1212	5	0
1	B	1227	0	1206	4	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	24	0	36	2	0
4	B	48	0	72	1	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	216	0	0	1	0
5	B	215	0	0	0	0
All	All	2949	0	2526	9	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 2.

All (9) 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:65[A]:LYS:NZ	5:A:312:HOH:O	2.39	0.54
1:A:19:CYS:HB2	4:A:202:EDO:H11	1.92	0.51
1:B:70:ARG:NE	4:B:202:EDO:O2	2.44	0.47
1:B:80[A]:ARG:HG2	1:B:80[A]:ARG:HH11	1.79	0.47
1:B:21:ASP:HB3	1:B:55:CYS:HB2	1.97	0.46
1:A:123:VAL:HG22	4:A:202:EDO:H12	2.01	0.43
1:A:107:GLU:HG2	1:A:110:ARG:NH2	2.34	0.43
1:B:24[B]:LEU:HD23	1:B:24[B]:LEU:C	2.41	0.41
1:A:107:GLU:HG2	1:A:110:ARG:HH22	1.86	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	155/153 (101%)	153 (99%)	2 (1%)	0	100	100
1	B	155/153 (101%)	153 (99%)	2 (1%)	0	100	100
All	All	310/306 (101%)	306 (99%)	4 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	130/131 (99%)	129 (99%)	1 (1%)	81	62
1	B	133/131 (102%)	133 (100%)	0	100	100
All	All	263/262 (100%)	262 (100%)	1 (0%)	91	80

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	HIS

Sometimes 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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 22 ligands modelled in this entry, 4 are monoatomic - leaving 18 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
4	EDO	B	212	-	3,3,3	0.56	0	2,2,2	0.06	0
4	EDO	B	208	-	3,3,3	0.28	0	2,2,2	0.51	0
4	EDO	A	204	-	3,3,3	0.42	0	2,2,2	0.42	0
4	EDO	B	204	-	3,3,3	0.41	0	2,2,2	0.22	0
4	EDO	B	209	-	3,3,3	0.47	0	2,2,2	0.32	0
4	EDO	B	202	-	3,3,3	0.42	0	2,2,2	0.35	0
4	EDO	B	205	-	3,3,3	0.47	0	2,2,2	0.37	0
4	EDO	B	211	-	3,3,3	0.42	0	2,2,2	0.36	0
4	EDO	A	203	-	3,3,3	0.42	0	2,2,2	0.52	0
4	EDO	A	202	-	3,3,3	0.42	0	2,2,2	0.25	0
4	EDO	B	203	-	3,3,3	0.51	0	2,2,2	0.20	0
4	EDO	A	206	-	3,3,3	0.53	0	2,2,2	0.43	0
4	EDO	B	206	-	3,3,3	0.51	0	2,2,2	0.15	0
4	EDO	A	205	-	3,3,3	0.42	0	2,2,2	0.25	0
4	EDO	B	207	-	3,3,3	0.54	0	2,2,2	0.14	0
4	EDO	A	207	-	3,3,3	0.48	0	2,2,2	0.30	0
4	EDO	B	210	-	3,3,3	0.55	0	2,2,2	0.16	0
4	EDO	B	213	-	3,3,3	0.62	0	2,2,2	0.18	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. '2' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	B	212	-	-	1/1/1/1	-
4	EDO	B	208	-	-	0/1/1/1	-
4	EDO	A	204	-	-	0/1/1/1	-
4	EDO	B	204	-	-	1/1/1/1	-
4	EDO	B	209	-	-	1/1/1/1	-
4	EDO	B	202	-	-	1/1/1/1	-
4	EDO	B	205	-	-	1/1/1/1	-
4	EDO	B	211	-	-	1/1/1/1	-
4	EDO	A	203	-	-	0/1/1/1	-
4	EDO	A	202	-	-	1/1/1/1	-
4	EDO	B	203	-	-	1/1/1/1	-
4	EDO	A	206	-	-	1/1/1/1	-
4	EDO	B	206	-	-	0/1/1/1	-
4	EDO	A	205	-	-	0/1/1/1	-

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	B	207	-	-	0/1/1/1	-
4	EDO	A	207	-	-	1/1/1/1	-
4	EDO	B	210	-	-	1/1/1/1	-
4	EDO	B	213	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	202	EDO	O1-C1-C2-O2
4	B	210	EDO	O1-C1-C2-O2
4	B	213	EDO	O1-C1-C2-O2
4	A	206	EDO	O1-C1-C2-O2
4	B	211	EDO	O1-C1-C2-O2
4	B	212	EDO	O1-C1-C2-O2
4	B	202	EDO	O1-C1-C2-O2
4	B	204	EDO	O1-C1-C2-O2
4	B	205	EDO	O1-C1-C2-O2
4	A	207	EDO	O1-C1-C2-O2
4	B	209	EDO	O1-C1-C2-O2
4	B	203	EDO	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	202	EDO	1	0
4	A	202	EDO	2	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

6.1 Protein, DNA and RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers

Unable to reproduce the depositors R factor - this section is therefore empty.