



wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 28, 2024 – 03:16 PM EST

PDB ID : 1EC7
Title : E. COLI GLUCARATE DEHYDRATASE NATIVE ENZYME
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Deposited on : 2000-01-25
Resolution : 1.90 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 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 : 2.36
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

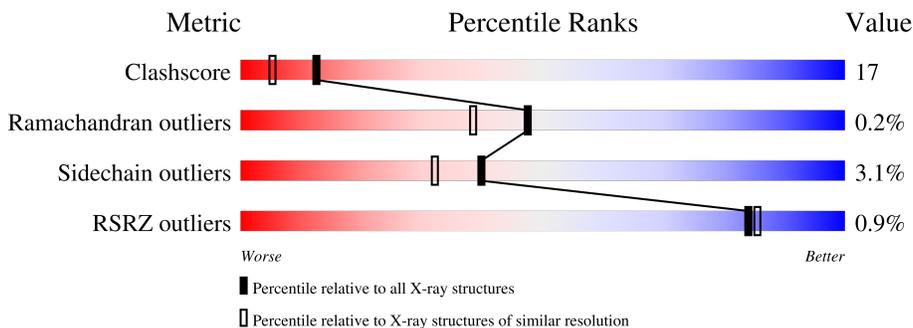
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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))
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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 $\leq 5\%$. 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	446	 2% 62% 29% 5% .
1	B	446	 2% 63% 31% 5% .
1	C	446	 67% 26% . .
1	D	446	 2% 60% 32% . .

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 14443 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.

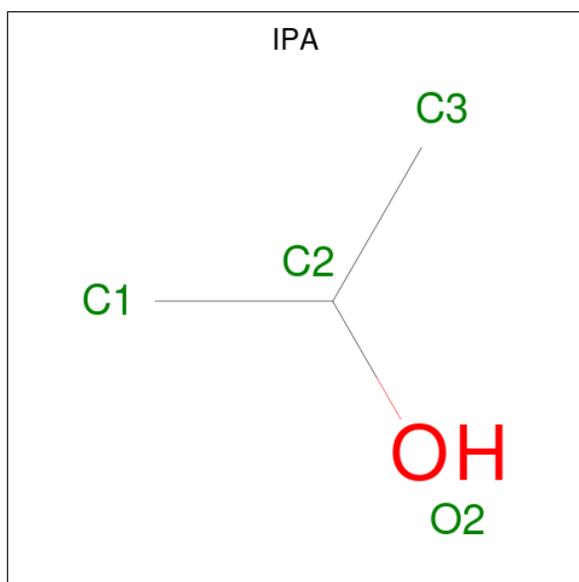
- Molecule 1 is a protein called GLUCARATE DEHYDRATASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	429	Total 3300	C 2091	N 571	O 617	S 21	0	0	0
1	B	440	Total 3398	C 2151	N 597	O 629	S 21	0	0	0
1	C	428	Total 3309	C 2096	N 576	O 616	S 21	0	1	0
1	D	431	Total 3314	C 2098	N 576	O 619	S 21	0	0	0

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total 1	Mg 1	0	0
2	B	1	Total 1	Mg 1	0	0
2	C	1	Total 1	Mg 1	0	0
2	D	1	Total 1	Mg 1	0	0

- Molecule 3 is ISOPROPYL ALCOHOL (three-letter code: IPA) (formula: C₃H₈O).



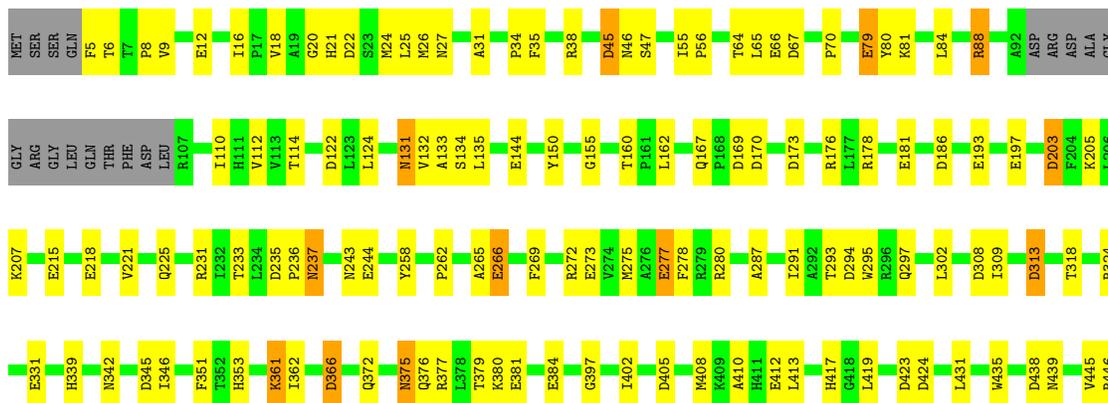
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 4 3 1	0	0
3	C	1	Total C O 4 3 1	0	0
3	D	1	Total C O 4 3 1	0	0
3	D	1	Total C O 4 3 1	0	0

- Molecule 4 is water.

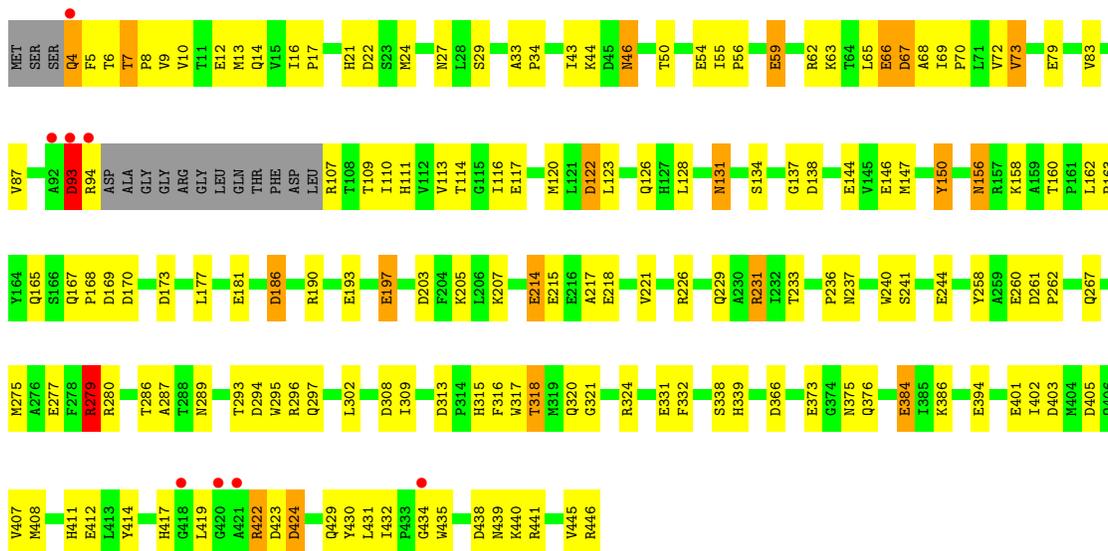
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	312	Total O 312 312	0	0
4	B	232	Total O 232 232	0	0
4	C	304	Total O 304 304	0	0
4	D	254	Total O 254 254	0	0



● Molecule 1: GLUCARATE DEHYDRATASE



● Molecule 1: GLUCARATE DEHYDRATASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	71.25Å 84.55Å 98.92Å 103.30° 94.00° 113.10°	Depositor
Resolution (Å)	30.00 – 1.90 19.86 – 1.70	Depositor EDS
% Data completeness (in resolution range)	93.0 (30.00-1.90) 87.7 (19.86-1.70)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.50 (at 1.70Å)	Xtrriage
Refinement program	TNT 5E	Depositor
R, R_{free}	0.194 , (Not available) 0.188 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	15.8	Xtrriage
Anisotropy	0.061	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 84.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	14443	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.18% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: IPA, MG

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	16/3377 (0.5%)	1.28	48/4582 (1.0%)
1	B	1.00	18/3476 (0.5%)	1.24	44/4710 (0.9%)
1	C	1.05	15/3386 (0.4%)	1.21	38/4592 (0.8%)
1	D	1.07	20/3391 (0.6%)	1.26	41/4602 (0.9%)
All	All	1.04	69/13630 (0.5%)	1.25	171/18486 (0.9%)

The worst 5 of 69 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	422	ARG	C-N	12.14	1.61	1.34
1	D	218	GLU	CD-OE2	8.67	1.35	1.25
1	A	79	GLU	CD-OE2	8.59	1.35	1.25
1	A	412	GLU	CD-OE2	8.37	1.34	1.25
1	D	144	GLU	CD-OE2	8.12	1.34	1.25

The worst 5 of 171 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	294	ASP	CB-CG-OD2	-10.04	109.26	118.30
1	A	366	ASP	CB-CG-OD1	8.84	126.26	118.30
1	A	173	ASP	CB-CG-OD2	-8.68	110.49	118.30
1	A	366	ASP	CB-CG-OD2	-8.48	110.67	118.30
1	B	22	ASP	CB-CG-OD2	-8.37	110.77	118.30

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	3300	0	3214	105	0
1	B	3398	0	3325	118	0
1	C	3309	0	3235	90	0
1	D	3314	0	3219	142	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	4	0	8	1	0
3	C	4	0	8	1	0
3	D	8	0	16	5	0
4	A	312	0	0	14	1
4	B	232	0	0	7	0
4	C	304	0	0	5	1
4	D	254	0	0	9	0
All	All	14443	0	13025	442	1

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 17.

The worst 5 of 442 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:D:429:GLN:NE2	1:D:435:TRP:H	1.43	1.15
1:D:21:HIS:H	1:D:376:GLN:HE22	1.05	0.96
1:B:414:TYR:HA	1:B:419:LEU:HD12	1.51	0.92
1:C:26:MET:HE2	1:C:31:ALA:HA	1.51	0.92
1:B:414:TYR:HA	1:B:419:LEU:CD1	2.00	0.91

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:2002:HOH:O	4:C:2036:HOH:O[1_455]	0.88	1.32

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	425/446 (95%)	409 (96%)	16 (4%)	0	100	100
1	B	436/446 (98%)	410 (94%)	25 (6%)	1 (0%)	47	38
1	C	425/446 (95%)	410 (96%)	14 (3%)	1 (0%)	47	38
1	D	427/446 (96%)	409 (96%)	17 (4%)	1 (0%)	47	38
All	All	1713/1784 (96%)	1638 (96%)	72 (4%)	3 (0%)	47	38

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	94	ARG
1	C	435	TRP
1	D	93	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	343/362 (95%)	331 (96%)	12 (4%)	36	27
1	B	353/362 (98%)	343 (97%)	10 (3%)	43	36
1	C	345/362 (95%)	336 (97%)	9 (3%)	46	39
1	D	343/362 (95%)	331 (96%)	12 (4%)	36	27
All	All	1384/1448 (96%)	1341 (97%)	43 (3%)	40	32

5 of 43 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	361	LYS
1	D	147	MET
1	C	375	ASN
1	D	7	THR
1	D	156	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 53 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	237	ASN
1	C	415	GLN
1	D	391	GLN
1	C	243	ASN
1	C	372	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](#)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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
3	IPA	D	602	-	3,3,3	0.50	0	3,3,3	0.30	0
3	IPA	D	604	-	3,3,3	0.50	0	3,3,3	0.38	0
3	IPA	A	603	-	3,3,3	0.68	0	3,3,3	0.18	0
3	IPA	C	601	-	3,3,3	0.55	0	3,3,3	0.27	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	602	IPA	2	0
3	D	604	IPA	3	0
3	A	603	IPA	1	0
3	C	601	IPA	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	D	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	D	422:ARG	C	423:ASP	N	1.62

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	429/446 (96%)	-0.45	4 (0%) 84 85	11, 23, 47, 88	0
1	B	440/446 (98%)	-0.31	3 (0%) 87 88	10, 30, 57, 81	0
1	C	428/446 (95%)	-0.56	0 100 100	10, 21, 45, 85	0
1	D	431/446 (96%)	-0.38	8 (1%) 66 69	11, 23, 57, 95	0
All	All	1728/1784 (96%)	-0.42	15 (0%) 84 85	10, 24, 53, 95	0

The worst 5 of 15 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	106	LEU	5.0
1	D	92	ALA	4.7
1	A	228	PRO	3.5
1	D	418	GLY	3.2
1	D	420	GLY	3.1

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](#)

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
3	IPA	C	601	4/4	0.91	0.12	23,26,30,35	0
3	IPA	D	604	4/4	0.91	0.12	21,29,30,41	0
3	IPA	A	603	4/4	0.92	0.14	17,21,53,54	0
3	IPA	D	602	4/4	0.93	0.12	16,30,40,44	0
2	MG	B	498	1/1	0.95	0.05	25,25,25,25	0
2	MG	C	498	1/1	0.96	0.07	24,24,24,24	0
2	MG	A	498	1/1	0.97	0.07	17,17,17,17	0
2	MG	D	498	1/1	0.97	0.05	23,23,23,23	0

6.5 Other polymers [i](#)

There are no such residues in this entry.