



wwPDB X-ray Structure Validation Summary Report ⓘ

May 15, 2020 – 02:44 am BST

PDB ID : 6HL8
Title : Crystal Structure of the CsiD Glutarate Hydroxylase in complex with Glutarate
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Deposited on : 2018-09-10
Resolution : 2.40 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 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	:	2.11
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

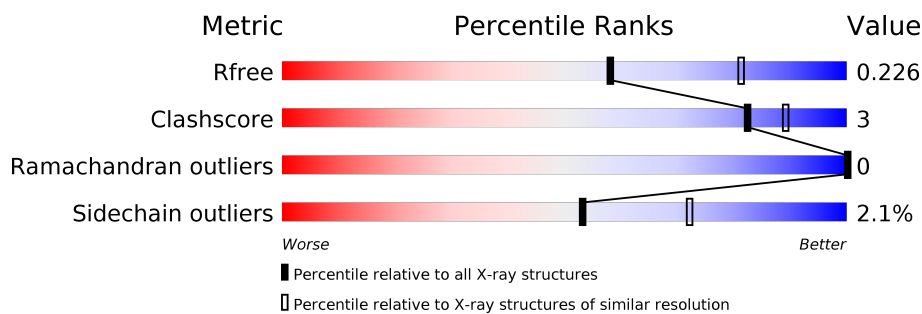
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)

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 $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	353	 75% 7% 18%
1	B	353	 76% 7% 17%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5111 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 Protein CsiD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	291	Total	C	N	O	S	25	2	0
			2397	1533	419	435	10			
1	B	292	Total	C	N	O	S	14	3	0
			2407	1541	421	435	10			

There are 56 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP P76621
A	-18	GLY	-	expression tag	UNP P76621
A	-17	SER	-	expression tag	UNP P76621
A	-16	SER	-	expression tag	UNP P76621
A	-15	HIS	-	expression tag	UNP P76621
A	-14	HIS	-	expression tag	UNP P76621
A	-13	HIS	-	expression tag	UNP P76621
A	-12	HIS	-	expression tag	UNP P76621
A	-11	HIS	-	expression tag	UNP P76621
A	-10	HIS	-	expression tag	UNP P76621
A	-9	SER	-	expression tag	UNP P76621
A	-8	SER	-	expression tag	UNP P76621
A	-7	GLY	-	expression tag	UNP P76621
A	-6	LEU	-	expression tag	UNP P76621
A	-5	VAL	-	expression tag	UNP P76621
A	-4	PRO	-	expression tag	UNP P76621
A	-3	ARG	-	expression tag	UNP P76621
A	-2	GLY	-	expression tag	UNP P76621
A	-1	SER	-	expression tag	UNP P76621
A	0	HIS	-	expression tag	UNP P76621
A	326	LEU	-	expression tag	UNP P76621
A	327	GLU	-	expression tag	UNP P76621
A	328	HIS	-	expression tag	UNP P76621
A	329	HIS	-	expression tag	UNP P76621
A	330	HIS	-	expression tag	UNP P76621

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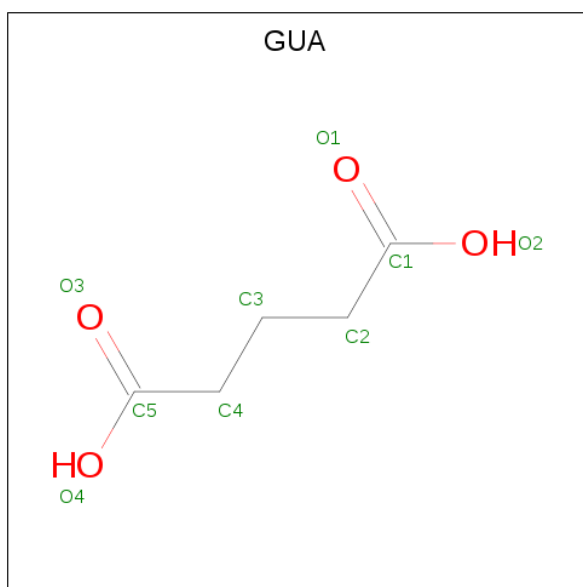
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Chain	Residue	Modelled	Actual	Comment	Reference
A	331	HIS	-	expression tag	UNP P76621
A	332	HIS	-	expression tag	UNP P76621
A	333	HIS	-	expression tag	UNP P76621
B	-19	MET	-	initiating methionine	UNP P76621
B	-18	GLY	-	expression tag	UNP P76621
B	-17	SER	-	expression tag	UNP P76621
B	-16	SER	-	expression tag	UNP P76621
B	-15	HIS	-	expression tag	UNP P76621
B	-14	HIS	-	expression tag	UNP P76621
B	-13	HIS	-	expression tag	UNP P76621
B	-12	HIS	-	expression tag	UNP P76621
B	-11	HIS	-	expression tag	UNP P76621
B	-10	HIS	-	expression tag	UNP P76621
B	-9	SER	-	expression tag	UNP P76621
B	-8	SER	-	expression tag	UNP P76621
B	-7	GLY	-	expression tag	UNP P76621
B	-6	LEU	-	expression tag	UNP P76621
B	-5	VAL	-	expression tag	UNP P76621
B	-4	PRO	-	expression tag	UNP P76621
B	-3	ARG	-	expression tag	UNP P76621
B	-2	GLY	-	expression tag	UNP P76621
B	-1	SER	-	expression tag	UNP P76621
B	0	HIS	-	expression tag	UNP P76621
B	326	LEU	-	expression tag	UNP P76621
B	327	GLU	-	expression tag	UNP P76621
B	328	HIS	-	expression tag	UNP P76621
B	329	HIS	-	expression tag	UNP P76621
B	330	HIS	-	expression tag	UNP P76621
B	331	HIS	-	expression tag	UNP P76621
B	332	HIS	-	expression tag	UNP P76621
B	333	HIS	-	expression tag	UNP P76621

- Molecule 2 is FE (II) ION (three-letter code: FE2) (formula: Fe).

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

- Molecule 3 is GLUTARIC ACID (three-letter code: GUA) (formula: C₅H₈O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			9	5	4		
3	B	1	Total	C	O	0	0
			9	5	4		

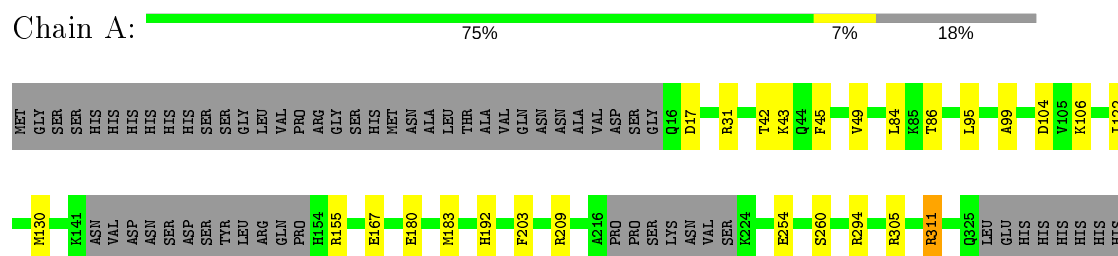
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	123	Total	O	0	0
			123	123		
4	B	164	Total	O	0	0
			164	164		

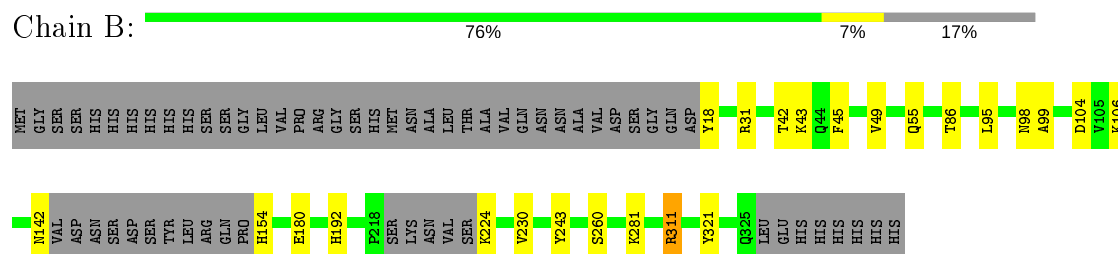
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: Protein CsiD



• Molecule 1: Protein CsiD



4 Data and refinement statistics

Property	Value	Source
Space group	P 4 21 2	Depositor
Cell constants a, b, c, α , β , γ	120.80 Å 120.80 Å 136.43 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.70 – 2.40 42.71 – 1.99	Depositor EDS
% Data completeness (in resolution range)	99.8 (29.70-2.40) 99.7 (42.71-1.99)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.27 (at 2.00 Å)	Xtriage
Refinement program	PHENIX (1.12_2829: ???)	Depositor
R, R_{free}	0.175 , 0.220 0.177 , 0.226	Depositor DCC
R_{free} test set	3530 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	29.3	Xtriage
Anisotropy	0.176	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 55.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5111	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 55.84 % 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 2.9854e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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: GUA, FE2

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	0.44	0/2460	0.58	0/3330
1	B	0.45	0/2475	0.60	0/3353
All	All	0.45	0/4935	0.59	0/6683

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	2397	0	2353	12	0
1	B	2407	0	2367	13	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	9	0	0	1	0
3	B	9	0	0	1	0
4	A	123	0	0	0	0
4	B	164	0	0	4	0
All	All	5111	0	4720	24	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 3.

The worst 5 of 24 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:104:ASP:OD1	1:A:106:LYS:HE3	1.90	0.72
1:A:311:ARG:NH1	3:A:402:GUA:O2	2.22	0.71
1:B:311:ARG:NH1	3:B:402:GUA:O1	2.34	0.60
1:B:224:LYS:N	4:B:503:HOH:O	2.35	0.58
1:A:155:ARG:O	1:A:294:ARG:NH2	2.43	0.52

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	287/353 (81%)	281 (98%)	6 (2%)	0	100	100
1	B	289/353 (82%)	283 (98%)	6 (2%)	0	100	100
All	All	576/706 (82%)	564 (98%)	12 (2%)	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	259/312 (83%)	253 (98%)	6 (2%)	50	70

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	261/312 (84%)	256 (98%)	5 (2%)	57	75
All	All	520/624 (83%)	509 (98%)	11 (2%)	53	72

5 of 11 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	260	SER
1	A	311	ARG
1	B	180	GLU
1	A	180	GLU
1	B	154	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 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	GUA	B	402	-	2,8,8	0.25	0	1,9,9	0.51	0
3	GUA	A	402	-	2,8,8	0.23	0	1,9,9	0.55	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. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GUA	B	402	-	-	1/2/6/6	-
3	GUA	A	402	-	-	0/2/6/6	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	402	GUA	C1-C2-C3-C4

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	402	GUA	1	0
3	A	402	GUA	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 ⓘ

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.