



wwPDB X-ray Structure Validation Summary Report ⓘ

May 16, 2020 – 11:01 am BST

PDB ID : 6GPN
Title : Crystal Structure of the CsiD Glutarate Hydroxylase in complex with N-Oxalylglycine
Authors : Williams, R.M.; Mayans, O.; Hartig, J.S.
Deposited on : 2018-06-06
Resolution : 2.20 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

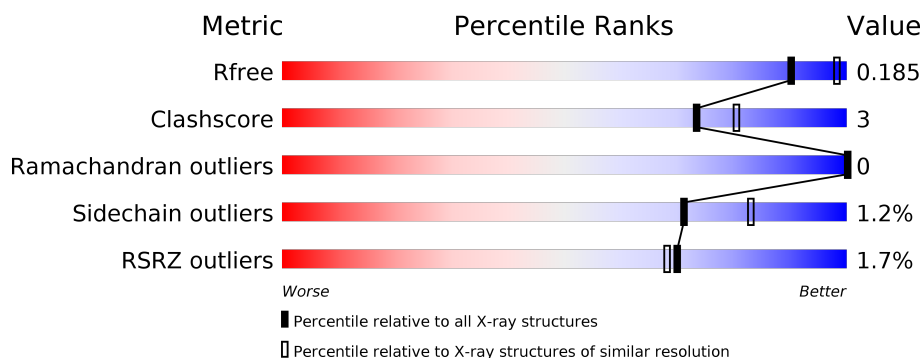
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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	353	<div> <div>2%</div> <div> <div></div> <div>75%</div> <div>7%</div> <div>18%</div> </div> </div>
1	B	353	<div> <div>%</div> <div> <div></div> <div>76%</div> <div>6%</div> <div>17%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4996 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	289	Total	C	N	O	S	0	1	0
			2336	1499	403	424	10			
1	B	292	Total	C	N	O	S	0	2	0
			2376	1524	411	431	10			

There are 56 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP A0A1X3JCQ2
A	-18	GLY	-	expression tag	UNP A0A1X3JCQ2
A	-17	SER	-	expression tag	UNP A0A1X3JCQ2
A	-16	SER	-	expression tag	UNP A0A1X3JCQ2
A	-15	HIS	-	expression tag	UNP A0A1X3JCQ2
A	-14	HIS	-	expression tag	UNP A0A1X3JCQ2
A	-13	HIS	-	expression tag	UNP A0A1X3JCQ2
A	-12	HIS	-	expression tag	UNP A0A1X3JCQ2
A	-11	HIS	-	expression tag	UNP A0A1X3JCQ2
A	-10	HIS	-	expression tag	UNP A0A1X3JCQ2
A	-9	SER	-	expression tag	UNP A0A1X3JCQ2
A	-8	SER	-	expression tag	UNP A0A1X3JCQ2
A	-7	GLY	-	expression tag	UNP A0A1X3JCQ2
A	-6	LEU	-	expression tag	UNP A0A1X3JCQ2
A	-5	VAL	-	expression tag	UNP A0A1X3JCQ2
A	-4	PRO	-	expression tag	UNP A0A1X3JCQ2
A	-3	ARG	-	expression tag	UNP A0A1X3JCQ2
A	-2	GLY	-	expression tag	UNP A0A1X3JCQ2
A	-1	SER	-	expression tag	UNP A0A1X3JCQ2
A	0	HIS	-	expression tag	UNP A0A1X3JCQ2
A	326	LEU	-	expression tag	UNP A0A1X3JCQ2
A	327	GLU	-	expression tag	UNP A0A1X3JCQ2
A	328	HIS	-	expression tag	UNP A0A1X3JCQ2
A	329	HIS	-	expression tag	UNP A0A1X3JCQ2
A	330	HIS	-	expression tag	UNP A0A1X3JCQ2

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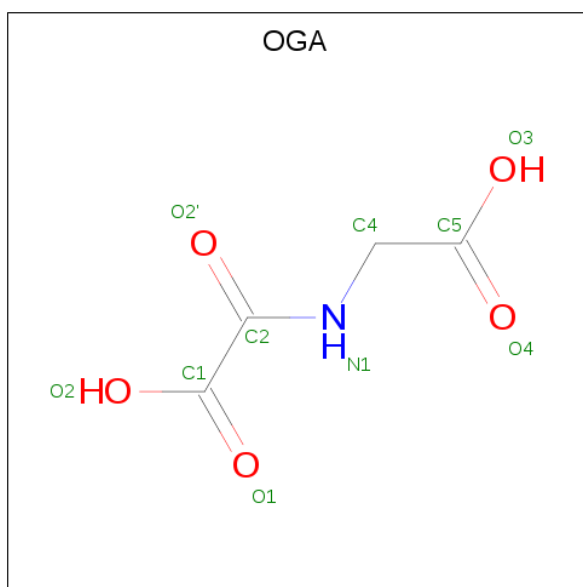
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Chain	Residue	Modelled	Actual	Comment	Reference
A	331	HIS	-	expression tag	UNP A0A1X3JCQ2
A	332	HIS	-	expression tag	UNP A0A1X3JCQ2
A	333	HIS	-	expression tag	UNP A0A1X3JCQ2
B	-19	MET	-	initiating methionine	UNP A0A1X3JCQ2
B	-18	GLY	-	expression tag	UNP A0A1X3JCQ2
B	-17	SER	-	expression tag	UNP A0A1X3JCQ2
B	-16	SER	-	expression tag	UNP A0A1X3JCQ2
B	-15	HIS	-	expression tag	UNP A0A1X3JCQ2
B	-14	HIS	-	expression tag	UNP A0A1X3JCQ2
B	-13	HIS	-	expression tag	UNP A0A1X3JCQ2
B	-12	HIS	-	expression tag	UNP A0A1X3JCQ2
B	-11	HIS	-	expression tag	UNP A0A1X3JCQ2
B	-10	HIS	-	expression tag	UNP A0A1X3JCQ2
B	-9	SER	-	expression tag	UNP A0A1X3JCQ2
B	-8	SER	-	expression tag	UNP A0A1X3JCQ2
B	-7	GLY	-	expression tag	UNP A0A1X3JCQ2
B	-6	LEU	-	expression tag	UNP A0A1X3JCQ2
B	-5	VAL	-	expression tag	UNP A0A1X3JCQ2
B	-4	PRO	-	expression tag	UNP A0A1X3JCQ2
B	-3	ARG	-	expression tag	UNP A0A1X3JCQ2
B	-2	GLY	-	expression tag	UNP A0A1X3JCQ2
B	-1	SER	-	expression tag	UNP A0A1X3JCQ2
B	0	HIS	-	expression tag	UNP A0A1X3JCQ2
B	326	LEU	-	expression tag	UNP A0A1X3JCQ2
B	327	GLU	-	expression tag	UNP A0A1X3JCQ2
B	328	HIS	-	expression tag	UNP A0A1X3JCQ2
B	329	HIS	-	expression tag	UNP A0A1X3JCQ2
B	330	HIS	-	expression tag	UNP A0A1X3JCQ2
B	331	HIS	-	expression tag	UNP A0A1X3JCQ2
B	332	HIS	-	expression tag	UNP A0A1X3JCQ2
B	333	HIS	-	expression tag	UNP A0A1X3JCQ2

- 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 N-OXALYLGLYCINE (three-letter code: OGA) (formula: C₄H₅NO₅).

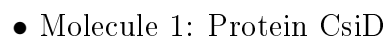


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	N	O	0	0
			10	4	1	5		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	100	Total	O	0	0
			100	100		
4	B	172	Total	O	0	0
			172	172		

- 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	121.38 Å 121.38 Å 136.57 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.91 – 2.20 42.91 – 1.99	Depositor EDS
% Data completeness (in resolution range)	99.9 (42.91-2.20) 99.8 (42.91-1.99)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.18	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.33 (at 1.98 Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
R, R_{free}	0.187 , 0.211 0.187 , 0.185	Depositor DCC
R_{free} test set	2180 reflections (3.10%)	wwPDB-VP
Wilson B-factor (Å ²)	37.1	Xtriage
Anisotropy	0.225	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 55.0	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	0.97	EDS
Total number of atoms	4996	wwPDB-VP
Average B, all atoms (Å ²)	52.0	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 28.91 % 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 1.6983e-03. 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: OGA, 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	0.41	1/2395 (0.0%)	0.60	1/3248 (0.0%)
1	B	0.42	0/2440	0.61	1/3308 (0.0%)
All	All	0.41	1/4835 (0.0%)	0.60	2/6556 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	40	GLN	CD-NE2	5.53	1.46	1.32

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	241	MET	CG-SD-CE	5.53	109.04	100.20
1	A	40	GLN	CB-CA-C	5.19	120.78	110.40

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	2336	0	2265	16	0
1	B	2376	0	2322	15	0
2	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1	0	0	0	0
3	B	10	0	3	2	0
4	A	100	0	0	1	0
4	B	172	0	0	2	0
All	All	4996	0	4590	31	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 31 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:180:GLU:OE2	1:B:183:MET:HG2	1.73	0.88
1:B:241:MET:HE3	1:B:291:LEU:HD11	1.58	0.84
1:A:182:ASN:HD22	1:A:302:ASP:HB2	1.51	0.74
1:B:157:MET:HE1	3:B:402:OGA:H4C1	1.76	0.67
1:B:141:LYS:NZ	4:B:503:HOH:O	2.32	0.61

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	284/353 (80%)	278 (98%)	6 (2%)	0	100	100
1	B	288/353 (82%)	283 (98%)	5 (2%)	0	100	100
All	All	572/706 (81%)	561 (98%)	11 (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	247/312 (79%)	245 (99%)	2 (1%)	81	90
1	B	255/312 (82%)	251 (98%)	4 (2%)	62	76
All	All	502/624 (80%)	496 (99%)	6 (1%)	71	83

5 of 6 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	209	ARG
1	B	322	GLN
1	B	260	SER
1	A	307	LEU
1	B	311	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	182	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry

Of 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 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	OGA	B	402	2	3,9,9	2.84	1 (33%)	4,11,11	3.24	2 (50%)

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	OGA	B	402	2	-	1/3/9/9	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	402	OGA	C2-N1	4.83	1.44	1.33

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	402	OGA	C1-C2-N1	4.87	120.44	115.60
3	B	402	OGA	C5-C4-N1	4.06	118.26	110.43

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	402	OGA	C5-C4-N1-C2

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	402	OGA	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 [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	289/353 (81%)	-0.21	6 (2%) 63 61	40, 52, 85, 117	0
1	B	292/353 (82%)	-0.23	4 (1%) 75 73	36, 45, 71, 103	0
All	All	581/706 (82%)	-0.22	10 (1%) 70 68	36, 48, 81, 117	0

The worst 5 of 10 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	17	ASP	3.8
1	B	218	PRO	3.7
1	A	40	GLN	2.6
1	A	138	PHE	2.5
1	B	224	LYS	2.5

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](#)

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	OGA	B	402	10/10	0.65	0.28	87,93,95,97	0
2	FE2	B	401	1/1	0.97	0.11	55,55,55,55	0
2	FE2	A	401	1/1	0.98	0.14	66,66,66,66	0

6.5 Other polymers [i](#)

There are no such residues in this entry.