



wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 31, 2016 – 08:02 PM GMT

PDB ID : 1IGW
Title : Crystal Structure of the Isocitrate Lyase from the A219C mutant of Escherichia coli
Authors : Britton, K.L.; Abeysinghe, I.S.B.; Baker, P.J.; Barynin, V.; Diehl, P.; Langridge, S.J.; McFadden, B.A.; Sedelnikova, S.E.; Stillman, T.J.; Weeradechapon, K.; Rice, D.W.
Deposited on : 2001-04-18
Resolution : 2.10 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
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

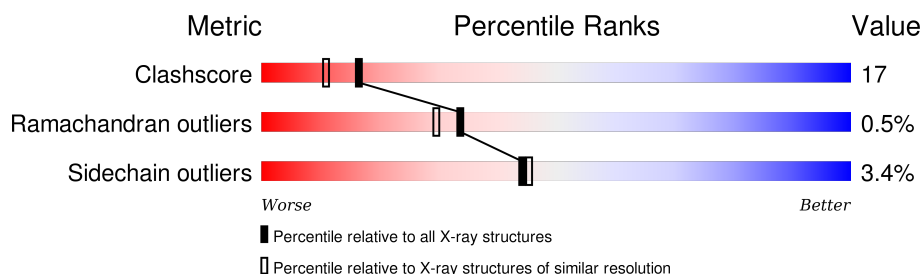
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)

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. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	434	
1	B	434	
1	C	434	
1	D	434	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	HG	C	438	-	-	X	-
2	HG	D	439	-	-	X	-
4	PYR	C	1446	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 13435 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 Isocitrate lyase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	396	Total	C	N	O	S	0	0	0
			3066	1948	521	583	14			
1	B	411	Total	C	N	O	S	0	0	0
			3191	2028	542	608	13			
1	C	416	Total	C	N	O	S	0	0	0
			3223	2046	549	613	15			
1	D	405	Total	C	N	O	S	0	0	0
			3139	1995	531	600	13			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	219	CYS	ALA	ENGINEERED	UNP P0A9G6
B	219	CYS	ALA	ENGINEERED	UNP P0A9G6
C	219	CYS	ALA	ENGINEERED	UNP P0A9G6
D	219	CYS	ALA	ENGINEERED	UNP P0A9G6

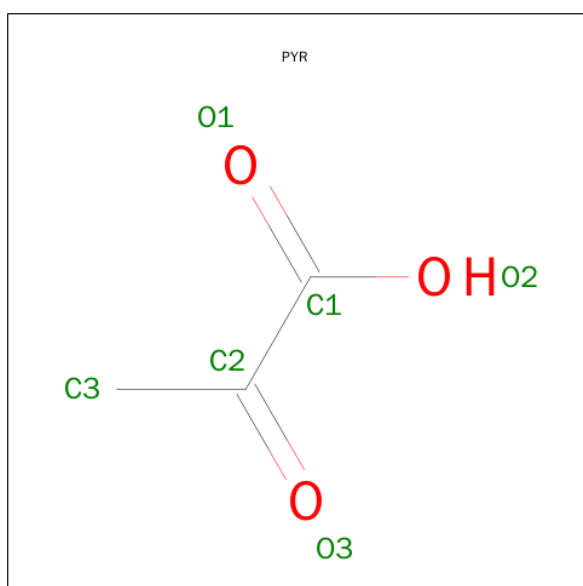
- Molecule 2 is MERCURY (II) ION (three-letter code: HG) (formula: Hg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	5	Total	Hg	0	0
			5	5		
2	A	5	Total	Hg	0	0
			5	5		
2	D	5	Total	Hg	0	0
			5	5		
2	C	6	Total	Hg	0	0
			6	6		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 4 is PYRUVIC ACID (three-letter code: PYR) (formula: $C_3H_4O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 6 3 3	0	0
4	B	1	Total C O 6 3 3	0	0
4	C	1	Total C O 6 3 3	0	0
4	D	1	Total C O 6 3 3	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	188	Total O 188 188	0	0
5	B	190	Total O 190 190	0	0

Continued on next page...

Continued from previous page...

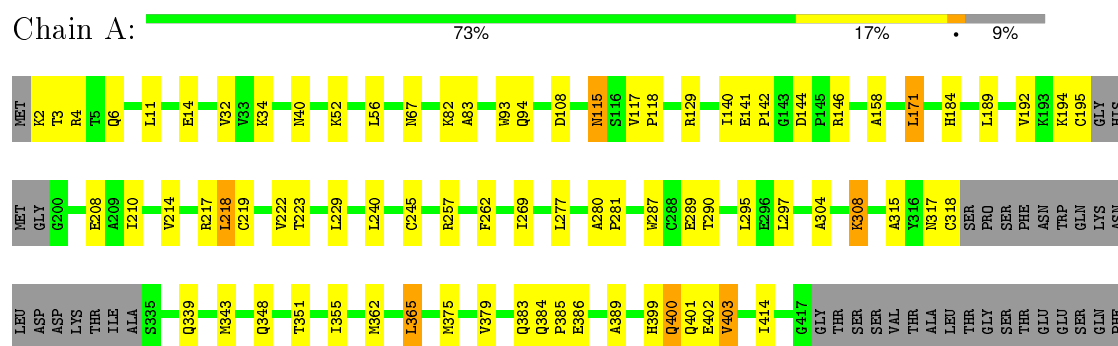
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	250	Total 250	O 250	0	0
5	D	139	Total 139	O 139	0	0

3 Residue-property plots

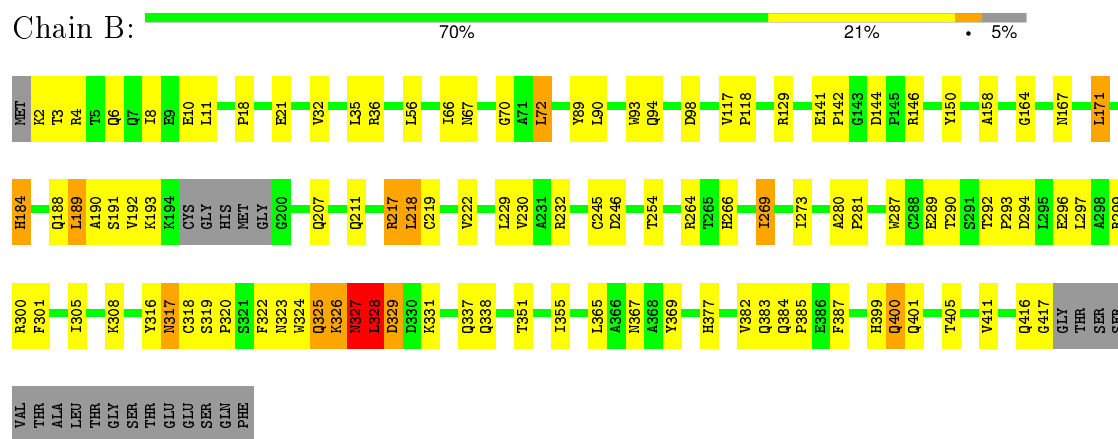
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

Note EDS was not executed.

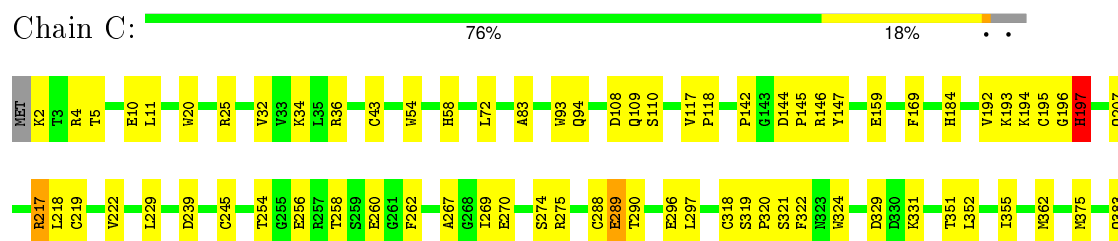
• Molecule 1: Isocitrate lyase

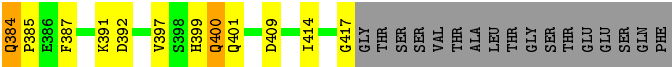


• Molecule 1: Isocitrate lyase

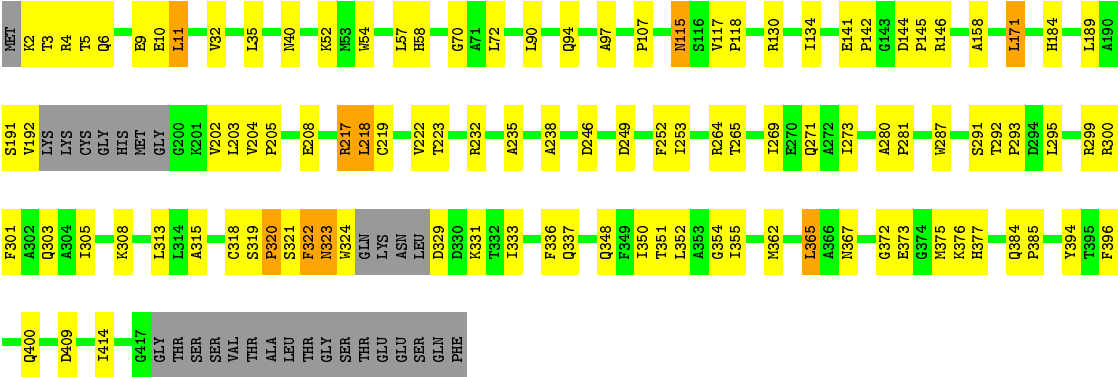


• Molecule 1: Isocitrate lyase





● Molecule 1: Isocitrate lyase



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	88.65Å 88.65Å 199.40Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	15.00 – 2.10	Depositor
% Data completeness (in resolution range)	(Not available) (15.00-2.10)	Depositor
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC	Depositor
R, R_{free}	0.184 , 0.235	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	13435	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

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: MG, PYR, HG

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	0.40	0/3134	0.94	4/4251 (0.1%)
1	B	0.44	1/3264 (0.0%)	1.06	17/4430 (0.4%)
1	C	0.41	0/3298	0.98	7/4476 (0.2%)
1	D	0.36	0/3211	0.88	3/4360 (0.1%)
All	All	0.40	1/12907 (0.0%)	0.97	31/17517 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	405	THR	C-N	-7.21	1.20	1.33

The worst 5 of 31 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	217	ARG	NE-CZ-NH1	9.95	125.27	120.30
1	A	2	LYS	N-CA-C	-9.24	86.04	111.00
1	B	72	LEU	CA-CB-CG	8.78	135.49	115.30
1	D	217	ARG	NE-CZ-NH2	-8.73	115.94	120.30
1	B	217	ARG	NE-CZ-NH2	-7.34	116.63	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	3066	0	2980	89	0
1	B	3191	0	3099	102	0
1	C	3223	0	3127	93	0
1	D	3139	0	3036	132	0
2	A	5	0	0	1	0
2	B	5	0	0	2	0
2	C	6	0	0	2	0
2	D	5	0	0	3	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	6	0	3	3	0
4	B	6	0	3	1	0
4	C	6	0	3	6	0
4	D	6	0	3	1	0
5	A	188	0	0	12	0
5	B	190	0	0	23	0
5	C	250	0	0	24	0
5	D	139	0	0	20	0
All	All	13435	0	12254	403	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 17.

The worst 5 of 403 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:293:PRO:HD3	1:D:322:PHE:CE1	1.53	1.43
1:D:293:PRO:HD3	1:D:322:PHE:CZ	1.56	1.39
1:D:409:ASP:HB3	5:D:1539:HOH:O	1.19	1.33
1:A:195:CYS:HB2	5:D:1565:HOH:O	1.30	1.25
1:B:416:GLN:OE1	1:C:195:CYS:SG	1.97	1.23

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	390/434 (90%)	375 (96%)	14 (4%)	1 (0%)	46	45
1	B	407/434 (94%)	390 (96%)	15 (4%)	2 (0%)	34	30
1	C	414/434 (95%)	401 (97%)	11 (3%)	2 (0%)	34	30
1	D	399/434 (92%)	381 (96%)	15 (4%)	3 (1%)	24	17
All	All	1610/1736 (93%)	1547 (96%)	55 (3%)	8 (0%)	34	30

5 of 8 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	326	LYS
1	C	197	HIS
1	D	322	PHE
1	D	323	ASN
1	B	328	LEU

5.3.2 Protein sidechains ⓘ

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	312/344 (91%)	302 (97%)	10 (3%)	46	48
1	B	326/344 (95%)	309 (95%)	17 (5%)	29	25
1	C	329/344 (96%)	320 (97%)	9 (3%)	52	56
1	D	320/344 (93%)	312 (98%)	8 (2%)	55	59
All	All	1287/1376 (94%)	1243 (97%)	44 (3%)	44	45

5 of 44 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	317	ASN
1	B	338	GLN
1	D	218	LEU
1	B	322	PHE
1	B	327	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 32 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	337	GLN
1	B	400	GLN
1	D	377	HIS
1	B	399	HIS
1	B	416	GLN

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 29 ligands modelled in this entry, 25 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
4	PYR	A	1444	3	2,5,5	3.99	1 (50%)	2,6,6	6.67	2 (100%)
4	PYR	B	1445	3	2,5,5	0.22	0	2,6,6	3.57	2 (100%)
4	PYR	C	1446	3	2,5,5	4.31	1 (50%)	2,6,6	1.19	0
4	PYR	D	1447	3	2,5,5	3.88	1 (50%)	2,6,6	3.69	2 (100%)

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
4	PYR	A	1444	3	-	0/0/4/4	0/0/0/0
4	PYR	B	1445	3	-	0/0/4/4	0/0/0/0
4	PYR	C	1446	3	-	0/0/4/4	0/0/0/0
4	PYR	D	1447	3	-	0/0/4/4	0/0/0/0

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	1447	PYR	O3-C2	5.42	1.41	1.22
4	A	1444	PYR	O3-C2	5.57	1.41	1.22
4	C	1446	PYR	O3-C2	6.09	1.43	1.22

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	1447	PYR	O3-C2-C3	-4.56	109.19	120.13
4	B	1445	PYR	O3-C2-C3	-4.48	109.41	120.13
4	A	1444	PYR	C3-C2-C1	-2.67	112.30	120.23
4	D	1447	PYR	C3-C2-C1	-2.54	112.67	120.23
4	B	1445	PYR	C3-C2-C1	-2.33	113.31	120.23

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1444	PYR	3	0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	1445	PYR	1	0
4	C	1446	PYR	6	0
4	D	1447	PYR	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

6.5 Other polymers [i](#)

EDS was not executed - this section will therefore be empty.