



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 14, 2017 – 02:08 AM EST

PDB ID : 5KGM
Title : 2.95Å resolution structure of Apo independent phosphoglycerate mutase from *C. elegans* (monoclinic form)
Authors : Lovell, S.; Mehzabeen, N.; Battaile, K.P.; Yu, H.; Dranchak, P.; MacArthur, R.; Li, Z.; Carlow, T.; Suga, H.; Inglese, J.
Deposited on : unknown
Resolution : 2.95 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

<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
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20030345
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20030345

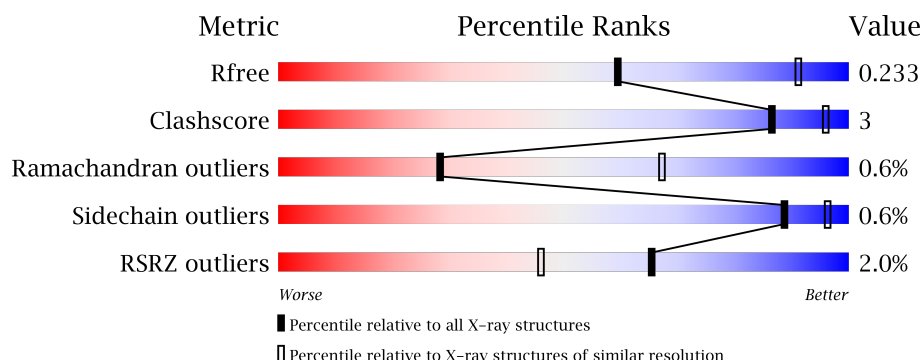
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 2.95 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}	100719	2395 (3.00-2.92)
Clashscore	112137	2773 (3.00-2.92)
Ramachandran outliers	110173	2680 (3.00-2.92)
Sidechain outliers	110143	2683 (3.00-2.92)
RSRZ outliers	101464	2421 (3.00-2.92)

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.

Mol	Chain	Length	Quality of chain
1	A	552	
1	B	552	

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	CL	A	602	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7845 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 2,3-bisphosphoglycerate-independent phosphoglycerate mutase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	519	Total	C	N	O	S	0	0	0
			3933	2483	681	741	28			
1	B	518	Total	C	N	O	S	0	0	0
			3904	2466	672	738	28			

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	540	LYS	-	expression tag	UNP G5EFZ1
A	541	LEU	-	expression tag	UNP G5EFZ1
A	542	ALA	-	expression tag	UNP G5EFZ1
A	543	ALA	-	expression tag	UNP G5EFZ1
A	544	ALA	-	expression tag	UNP G5EFZ1
A	545	LEU	-	expression tag	UNP G5EFZ1
A	546	GLU	-	expression tag	UNP G5EFZ1
A	547	HIS	-	expression tag	UNP G5EFZ1
A	548	HIS	-	expression tag	UNP G5EFZ1
A	549	HIS	-	expression tag	UNP G5EFZ1
A	550	HIS	-	expression tag	UNP G5EFZ1
A	551	HIS	-	expression tag	UNP G5EFZ1
A	552	HIS	-	expression tag	UNP G5EFZ1
B	540	LYS	-	expression tag	UNP G5EFZ1
B	541	LEU	-	expression tag	UNP G5EFZ1
B	542	ALA	-	expression tag	UNP G5EFZ1
B	543	ALA	-	expression tag	UNP G5EFZ1
B	544	ALA	-	expression tag	UNP G5EFZ1
B	545	LEU	-	expression tag	UNP G5EFZ1
B	546	GLU	-	expression tag	UNP G5EFZ1
B	547	HIS	-	expression tag	UNP G5EFZ1
B	548	HIS	-	expression tag	UNP G5EFZ1
B	549	HIS	-	expression tag	UNP G5EFZ1
B	550	HIS	-	expression tag	UNP G5EFZ1
B	551	HIS	-	expression tag	UNP G5EFZ1

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Chain	Residue	Modelled	Actual	Comment	Reference
B	552	HIS	-	expression tag	UNP G5EFZ1

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	2	Total Cl 2 2	0	0
2	A	2	Total Cl 2 2	0	0

- Molecule 3 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

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

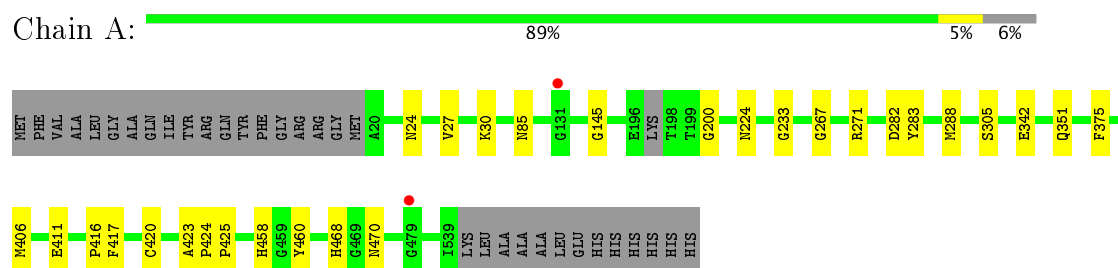
- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

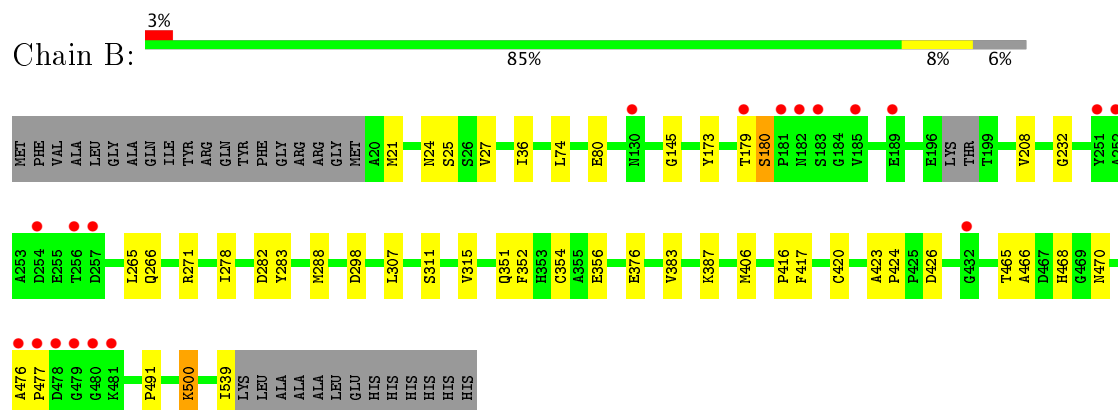
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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.

- Molecule 1: 2,3-bisphosphoglycerate-independent phosphoglycerate mutase



- Molecule 1: 2,3-bisphosphoglycerate-independent phosphoglycerate mutase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	67.85Å 94.38Å 102.20Å 90.00° 96.57° 90.00°	Depositor
Resolution (Å)	42.98 – 2.95 46.48 – 2.95	Depositor EDS
% Data completeness (in resolution range)	99.6 (42.98-2.95) 99.6 (46.48-2.95)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.24 (at 2.96Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.184 , 0.237 0.187 , 0.233	Depositor DCC
R_{free} test set	1352 reflections (5.01%)	DCC
Wilson B-factor (Å ²)	48.4	Xtriage
Anisotropy	0.472	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 43.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7845	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.23% 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: ZN, MN, CL

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	0.39	0/4021	0.62	0/5459
1	B	0.39	0/3992	0.62	1/5424 (0.0%)
All	All	0.39	0/8013	0.62	1/10883 (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	426	ASP	CB-CG-OD1	7.50	125.05	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	3933	0	3776	15	0
1	B	3904	0	3726	25	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	7845	0	7502	39	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.

All (39) 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:465:THR:HG22	1:B:466:ALA:H	1.58	0.68
1:B:476:ALA:HB1	1:B:477:PRO:HD2	1.88	0.55
1:A:233:GLY:N	1:A:271:ARG:HD2	2.21	0.55
1:A:145:GLY:O	1:A:283:TYR:OH	2.30	0.49
1:A:24:ASN:O	1:A:27:VAL:HG12	2.12	0.49
1:A:423:ALA:N	1:A:424:PRO:CD	2.76	0.48
1:B:465:THR:HG22	1:B:466:ALA:N	2.27	0.48
1:A:351:GLN:HA	1:A:417:PHE:O	2.14	0.48
1:B:74:LEU:HD11	1:B:80:GLU:HB2	1.95	0.48
1:B:24:ASN:O	1:B:27:VAL:HG12	2.14	0.47
1:B:356:GLU:HA	1:B:383:VAL:O	2.14	0.46
1:B:21:MET:SD	1:B:376:GLU:HG2	2.57	0.45
1:B:145:GLY:O	1:B:283:TYR:OH	2.34	0.45
1:B:352:PHE:CE2	1:B:354:CYS:HB2	2.51	0.45
1:A:85:ASN:ND2	1:B:387:LYS:HG2	2.31	0.45
1:A:267:GLY:O	1:A:271:ARG:NH1	2.50	0.45
1:B:36:ILE:O	1:B:466:ALA:HA	2.17	0.45
1:A:30:LYS:HB3	1:A:460:TYR:CE1	2.52	0.43
1:B:423:ALA:N	1:B:424:PRO:CD	2.81	0.43
1:B:179:THR:O	1:B:179:THR:HG23	2.19	0.43
1:B:179:THR:O	1:B:180:SER:O	2.37	0.42
1:B:265:LEU:O	1:B:266:GLN:HB2	2.19	0.42
1:B:278:ILE:CG2	1:B:315:VAL:HG22	2.49	0.42
1:B:232:GLY:HA3	1:B:307:LEU:HD13	2.00	0.42
1:B:232:GLY:HA3	1:B:307:LEU:CD1	2.50	0.41
1:B:465:THR:HG23	1:B:491:PRO:O	2.20	0.41
1:B:173:TYR:HA	1:B:208:VAL:O	2.21	0.41
1:B:282:ASP:HB2	1:B:288:MET:HE3	2.01	0.41
1:B:351:GLN:HA	1:B:417:PHE:O	2.21	0.41
1:A:406:MET:SD	1:A:420:CYS:HB3	2.61	0.41
1:A:27:VAL:HG11	1:A:416:PRO:HB3	2.03	0.41
1:A:411:GLU:HG2	1:A:458:HIS:NE2	2.36	0.41
1:B:406:MET:SD	1:B:420:CYS:HB3	2.60	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:282:ASP:HB2	1:A:288:MET:HE3	2.02	0.41
1:A:342:GLU:HB2	1:A:375:PHE:CZ	2.56	0.41
1:A:224:ASN:OD1	1:A:305:SER:HB2	2.20	0.40
1:B:27:VAL:HG11	1:B:416:PRO:HB3	2.03	0.40
1:A:424:PRO:HB2	1:A:425:PRO:HD3	2.03	0.40
1:B:500:LYS:HE2	1:B:539:ILE:HG21	2.04	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	515/552 (93%)	504 (98%)	8 (2%)	3 (1%)	28	67
1	B	514/552 (93%)	500 (97%)	11 (2%)	3 (1%)	28	67
All	All	1029/1104 (93%)	1004 (98%)	19 (2%)	6 (1%)	28	67

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	180	SER
1	A	470	ASN
1	B	470	ASN
1	B	468	HIS
1	A	200	GLY
1	A	468	HIS

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	403/445 (91%)	403 (100%)	0	100	100
1	B	398/445 (89%)	393 (99%)	5 (1%)	73	91
All	All	801/890 (90%)	796 (99%)	5 (1%)	89	96

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	25	SER
1	B	271	ARG
1	B	298	ASP
1	B	311	SER
1	B	500	LYS

Some 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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 8 are monoatomic - leaving 0 for Mogul analysis.

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.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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	519/552 (94%)	-0.05	2 (0%) 92 82	29, 45, 60, 95	0
1	B	518/552 (93%)	0.13	19 (3%) 42 27	33, 50, 75, 98	0
All	All	1037/1104 (93%)	0.04	21 (2%) 65 46	29, 47, 68, 98	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	252	ALA	3.7
1	B	476	ALA	3.3
1	B	185	VAL	3.3
1	B	182	ASN	3.2
1	B	183	SER	3.1
1	B	251	TYR	3.0
1	B	179	THR	2.7
1	B	479	GLY	2.7
1	B	254	ASP	2.6
1	B	130	ASN	2.5
1	B	256	THR	2.5
1	B	481	LYS	2.5
1	B	432	GLY	2.4
1	B	477	PRO	2.4
1	B	257	ASP	2.4
1	A	479	GLY	2.3
1	B	480	GLY	2.2
1	B	478	ASP	2.2
1	B	189	GLU	2.1
1	B	181	PRO	2.1
1	A	131	GLY	2.0

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	CL	A	602	1/1	0.91	0.28	3.23	44,44,44,44	0
3	MN	A	603	1/1	0.95	0.20	0.78	52,52,52,52	0
3	MN	B	603	1/1	0.94	0.18	0.09	45,45,45,45	0
4	ZN	A	604	1/1	0.99	0.18	-0.46	54,54,54,54	0
4	ZN	B	604	1/1	1.00	0.08	-4.45	41,41,41,41	0
2	CL	B	602	1/1	0.98	0.37	-	49,49,49,49	0
2	CL	B	601	1/1	0.96	0.20	-	37,37,37,37	0
2	CL	A	601	1/1	0.96	0.21	-	34,34,34,34	0

6.5 Other polymers [i](#)

There are no such residues in this entry.