



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 04:25 AM UTC

PDB ID : 1YJX / pdb_00001yjsx
Title : Crystal structure of human B type phosphoglycerate mutase
Authors : Wang, Y.; Wei, Z.; Liu, L.; Gong, W.
Deposited on : 2005-01-16
Resolution : 2.80 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	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

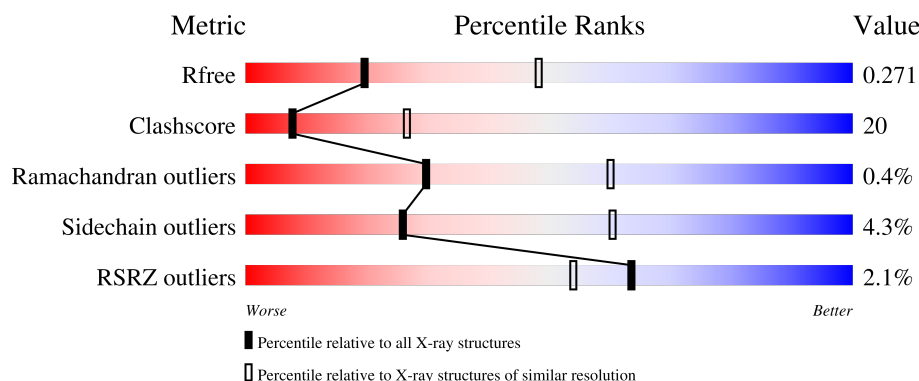
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	262	2% 61% 29% • 6%
1	B	262	% 59% 31% • 8%
1	C	262	2% 58% 32% • 7%
1	D	262	% 60% 32% • 6%
1	E	262	2% 62% 30% • 7%

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Mol	Chain	Length	Quality of chain
1	F	262	
1	G	262	
1	H	262	
1	I	262	
1	J	262	
1	K	262	
1	L	262	

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
3	CIT	B	602	-	X	-	-
3	CIT	C	603	-	X	-	-
3	CIT	D	604	-	X	-	-
3	CIT	E	605	-	X	-	-
3	CIT	F	606	-	X	-	-
3	CIT	H	608	-	X	-	-
3	CIT	I	609	-	X	-	-
3	CIT	J	610	-	X	-	-
3	CIT	K	611	-	X	-	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 23286 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 Phosphoglycerate mutase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	245	Total	C	N	O	S	0	0	0
			1913	1218	337	351	7			
1	B	242	Total	C	N	O	S	0	0	0
			1903	1211	335	350	7			
1	C	243	Total	C	N	O	S	0	0	0
			1898	1209	332	351	6			
1	D	245	Total	C	N	O	S	0	0	0
			1914	1218	336	353	7			
1	E	243	Total	C	N	O	S	0	0	0
			1897	1208	332	351	6			
1	F	242	Total	C	N	O	S	0	0	0
			1890	1204	332	347	7			
1	G	245	Total	C	N	O	S	0	0	0
			1887	1200	329	351	7			
1	H	241	Total	C	N	O	S	0	0	0
			1875	1196	328	345	6			
1	I	242	Total	C	N	O	S	0	0	0
			1904	1213	332	352	7			
1	J	244	Total	C	N	O	S	0	0	0
			1918	1220	339	352	7			
1	K	239	Total	C	N	O	S	0	0	0
			1870	1193	332	340	5			
1	L	245	Total	C	N	O	S	0	0	0
			1921	1222	342	350	7			

There are 96 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	255	LEU	-	expression tag	UNP P18669
A	256	GLU	-	expression tag	UNP P18669
A	257	HIS	-	expression tag	UNP P18669
A	258	HIS	-	expression tag	UNP P18669
A	259	HIS	-	expression tag	UNP P18669

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Chain	Residue	Modelled	Actual	Comment	Reference
A	260	HIS	-	expression tag	UNP P18669
A	261	HIS	-	expression tag	UNP P18669
A	262	HIS	-	expression tag	UNP P18669
B	255	LEU	-	expression tag	UNP P18669
B	256	GLU	-	expression tag	UNP P18669
B	257	HIS	-	expression tag	UNP P18669
B	258	HIS	-	expression tag	UNP P18669
B	259	HIS	-	expression tag	UNP P18669
B	260	HIS	-	expression tag	UNP P18669
B	261	HIS	-	expression tag	UNP P18669
B	262	HIS	-	expression tag	UNP P18669
C	255	LEU	-	expression tag	UNP P18669
C	256	GLU	-	expression tag	UNP P18669
C	257	HIS	-	expression tag	UNP P18669
C	258	HIS	-	expression tag	UNP P18669
C	259	HIS	-	expression tag	UNP P18669
C	260	HIS	-	expression tag	UNP P18669
C	261	HIS	-	expression tag	UNP P18669
C	262	HIS	-	expression tag	UNP P18669
D	255	LEU	-	expression tag	UNP P18669
D	256	GLU	-	expression tag	UNP P18669
D	257	HIS	-	expression tag	UNP P18669
D	258	HIS	-	expression tag	UNP P18669
D	259	HIS	-	expression tag	UNP P18669
D	260	HIS	-	expression tag	UNP P18669
D	261	HIS	-	expression tag	UNP P18669
D	262	HIS	-	expression tag	UNP P18669
E	255	LEU	-	expression tag	UNP P18669
E	256	GLU	-	expression tag	UNP P18669
E	257	HIS	-	expression tag	UNP P18669
E	258	HIS	-	expression tag	UNP P18669
E	259	HIS	-	expression tag	UNP P18669
E	260	HIS	-	expression tag	UNP P18669
E	261	HIS	-	expression tag	UNP P18669
E	262	HIS	-	expression tag	UNP P18669
F	255	LEU	-	expression tag	UNP P18669
F	256	GLU	-	expression tag	UNP P18669
F	257	HIS	-	expression tag	UNP P18669
F	258	HIS	-	expression tag	UNP P18669
F	259	HIS	-	expression tag	UNP P18669
F	260	HIS	-	expression tag	UNP P18669
F	261	HIS	-	expression tag	UNP P18669

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Chain	Residue	Modelled	Actual	Comment	Reference
F	262	HIS	-	expression tag	UNP P18669
G	255	LEU	-	expression tag	UNP P18669
G	256	GLU	-	expression tag	UNP P18669
G	257	HIS	-	expression tag	UNP P18669
G	258	HIS	-	expression tag	UNP P18669
G	259	HIS	-	expression tag	UNP P18669
G	260	HIS	-	expression tag	UNP P18669
G	261	HIS	-	expression tag	UNP P18669
G	262	HIS	-	expression tag	UNP P18669
H	255	LEU	-	expression tag	UNP P18669
H	256	GLU	-	expression tag	UNP P18669
H	257	HIS	-	expression tag	UNP P18669
H	258	HIS	-	expression tag	UNP P18669
H	259	HIS	-	expression tag	UNP P18669
H	260	HIS	-	expression tag	UNP P18669
H	261	HIS	-	expression tag	UNP P18669
H	262	HIS	-	expression tag	UNP P18669
I	255	LEU	-	expression tag	UNP P18669
I	256	GLU	-	expression tag	UNP P18669
I	257	HIS	-	expression tag	UNP P18669
I	258	HIS	-	expression tag	UNP P18669
I	259	HIS	-	expression tag	UNP P18669
I	260	HIS	-	expression tag	UNP P18669
I	261	HIS	-	expression tag	UNP P18669
I	262	HIS	-	expression tag	UNP P18669
J	255	LEU	-	expression tag	UNP P18669
J	256	GLU	-	expression tag	UNP P18669
J	257	HIS	-	expression tag	UNP P18669
J	258	HIS	-	expression tag	UNP P18669
J	259	HIS	-	expression tag	UNP P18669
J	260	HIS	-	expression tag	UNP P18669
J	261	HIS	-	expression tag	UNP P18669
J	262	HIS	-	expression tag	UNP P18669
K	255	LEU	-	expression tag	UNP P18669
K	256	GLU	-	expression tag	UNP P18669
K	257	HIS	-	expression tag	UNP P18669
K	258	HIS	-	expression tag	UNP P18669
K	259	HIS	-	expression tag	UNP P18669
K	260	HIS	-	expression tag	UNP P18669
K	261	HIS	-	expression tag	UNP P18669
K	262	HIS	-	expression tag	UNP P18669
L	255	LEU	-	expression tag	UNP P18669

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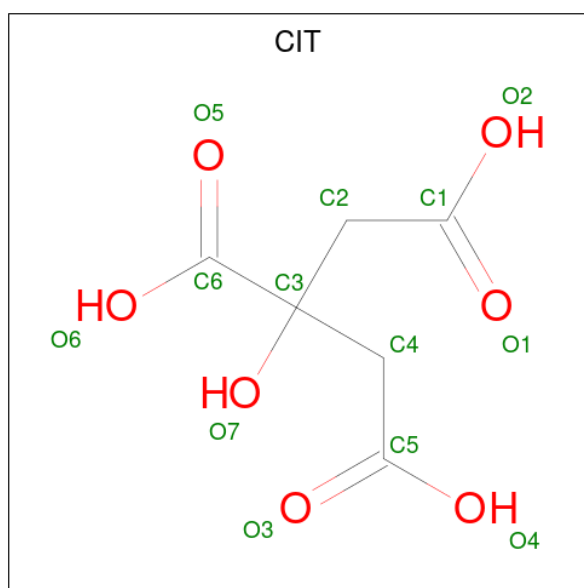
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Chain	Residue	Modelled	Actual	Comment	Reference
L	256	GLU	-	expression tag	UNP P18669
L	257	HIS	-	expression tag	UNP P18669
L	258	HIS	-	expression tag	UNP P18669
L	259	HIS	-	expression tag	UNP P18669
L	260	HIS	-	expression tag	UNP P18669
L	261	HIS	-	expression tag	UNP P18669
L	262	HIS	-	expression tag	UNP P18669

- Molecule 2 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Cl 1 1	0	0
2	C	1	Total Cl 1 1	0	0
2	F	1	Total Cl 1 1	0	0
2	H	1	Total Cl 1 1	0	0
2	J	1	Total Cl 1 1	0	0
2	K	1	Total Cl 1 1	0	0

- Molecule 3 is CITRIC ACID (CCD ID: CIT) (formula: C₆H₈O₇).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 13 6 7	0	0
3	B	1	Total C O 13 6 7	0	0
3	C	1	Total C O 13 6 7	0	0
3	D	1	Total C O 13 6 7	0	0
3	E	1	Total C O 13 6 7	0	0
3	F	1	Total C O 13 6 7	0	0
3	G	1	Total C O 13 6 7	0	0
3	H	1	Total C O 13 6 7	0	0
3	I	1	Total C O 13 6 7	0	0
3	J	1	Total C O 13 6 7	0	0
3	K	1	Total C O 13 6 7	0	0
3	L	1	Total C O 13 6 7	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	43	Total O 43 43	0	0
4	B	45	Total O 45 45	0	0
4	C	21	Total O 21 21	0	0
4	D	33	Total O 33 33	0	0
4	E	17	Total O 17 17	0	0
4	F	27	Total O 27 27	0	0
4	G	12	Total O 12 12	0	0
4	H	18	Total O 18 18	0	0

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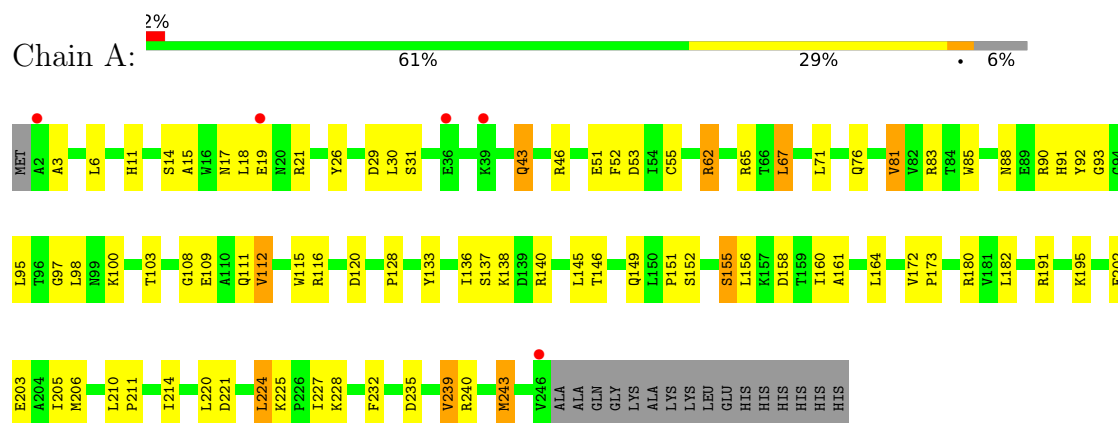
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	I	19	Total 19	O 19	0	0
4	J	15	Total 15	O 15	0	0
4	K	35	Total 35	O 35	0	0
4	L	49	Total 49	O 49	0	0

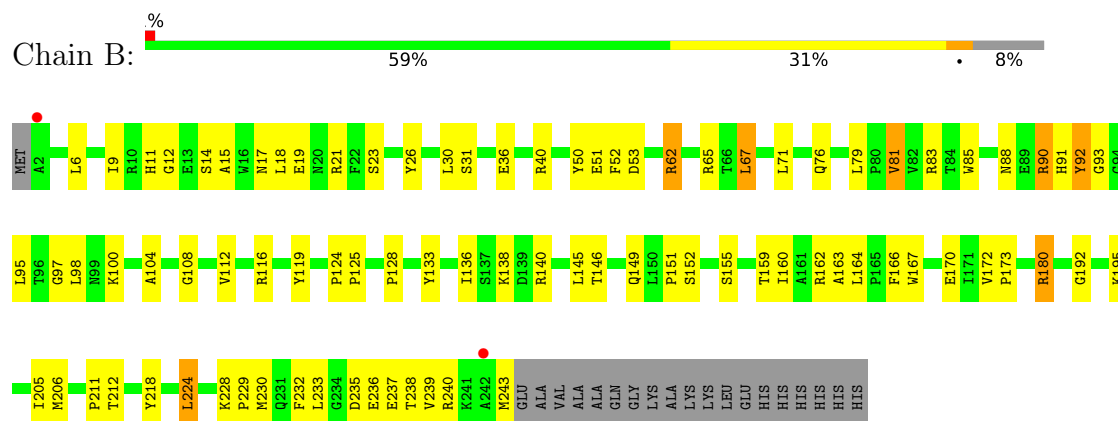
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

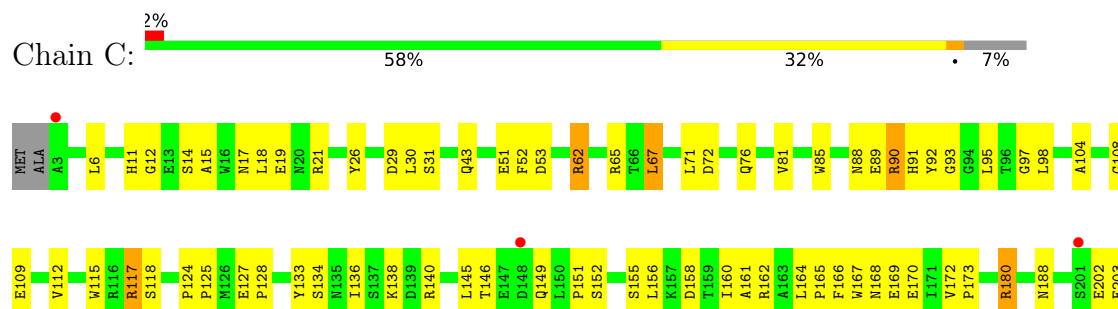
• Molecule 1: Phosphoglycerate mutase 1



• Molecule 1: Phosphoglycerate mutase 1

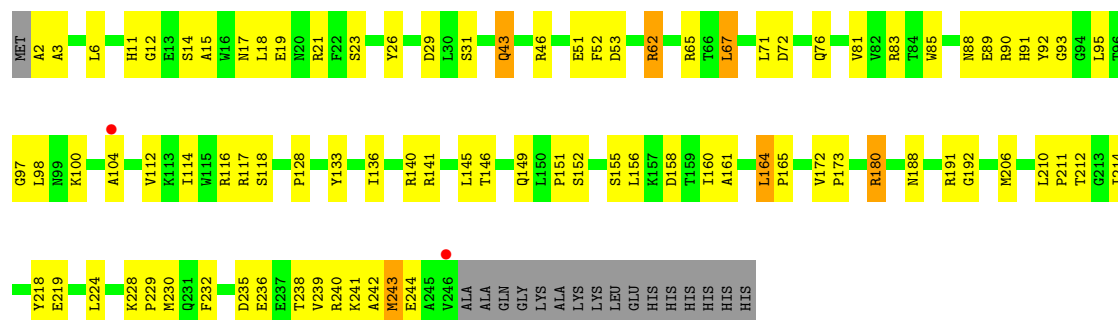


• Molecule 1: Phosphoglycerate mutase 1

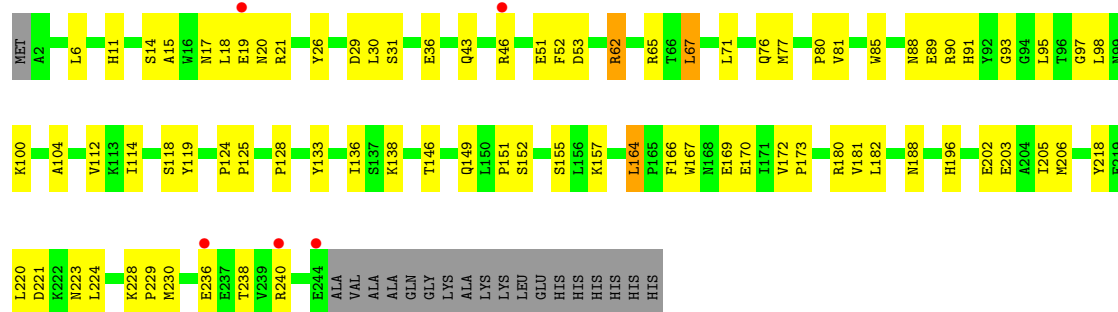




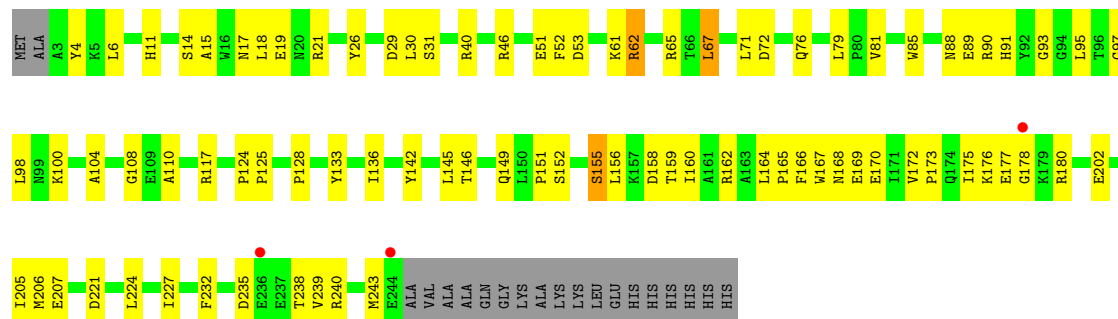
• Molecule 1: Phosphoglycerate mutase 1



• Molecule 1: Phosphoglycerate mutase 1

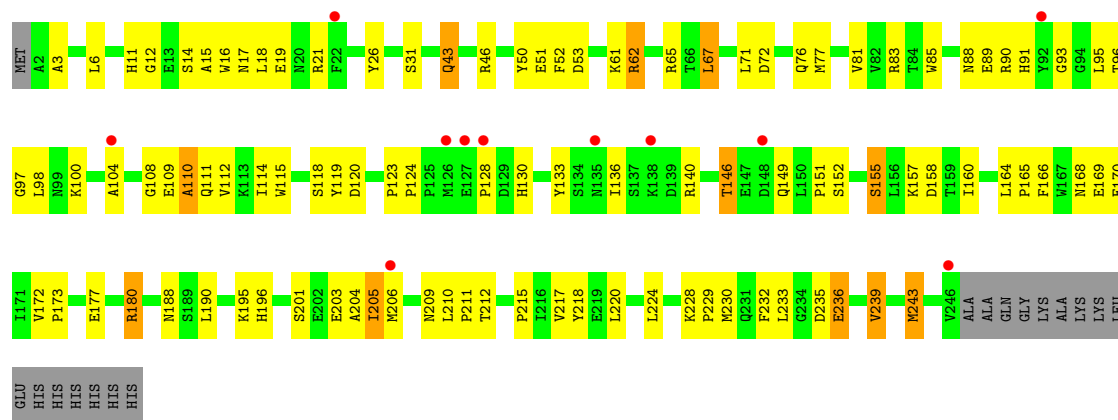


• Molecule 1: Phosphoglycerate mutase 1

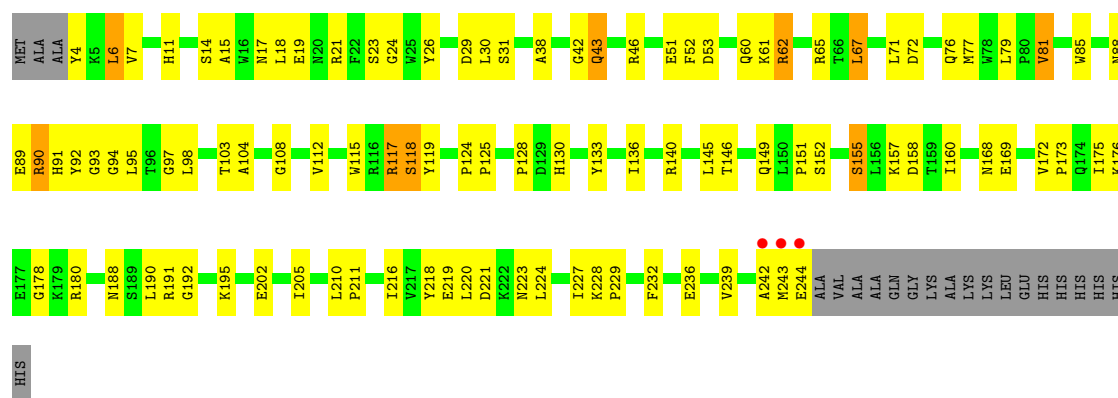


• Molecule 1: Phosphoglycerate mutase 1

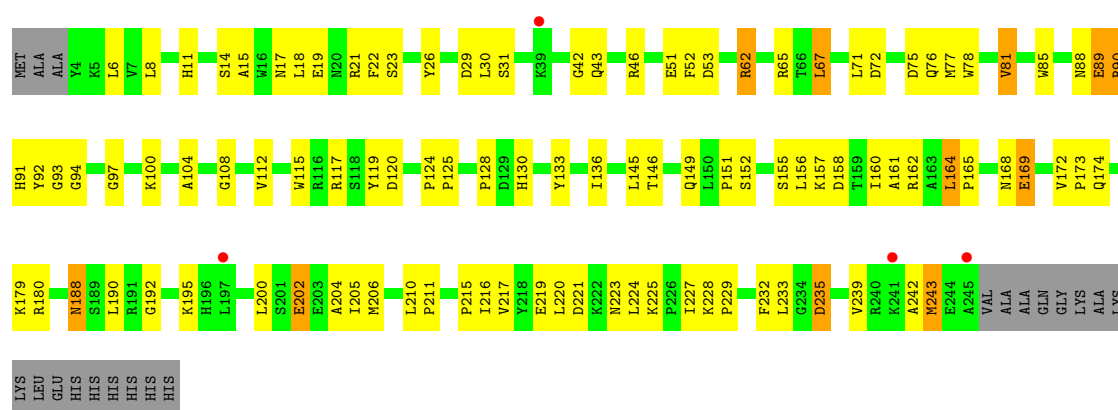




• Molecule 1: Phosphoglycerate mutase 1

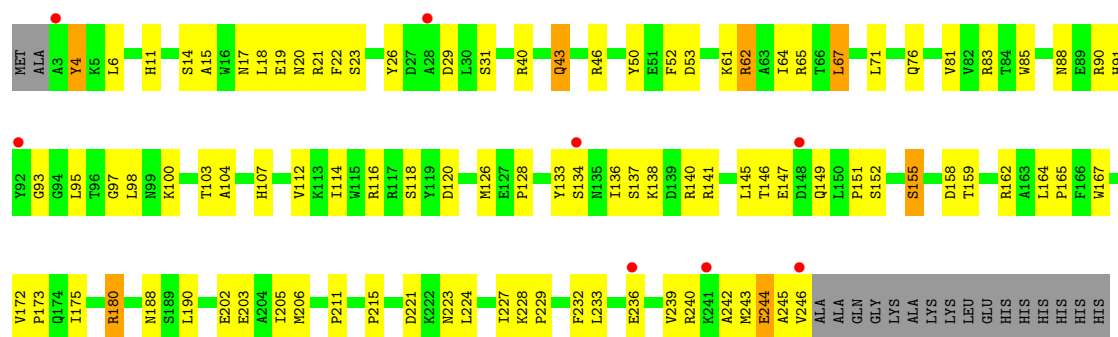


• Molecule 1: Phosphoglycerate mutase 1

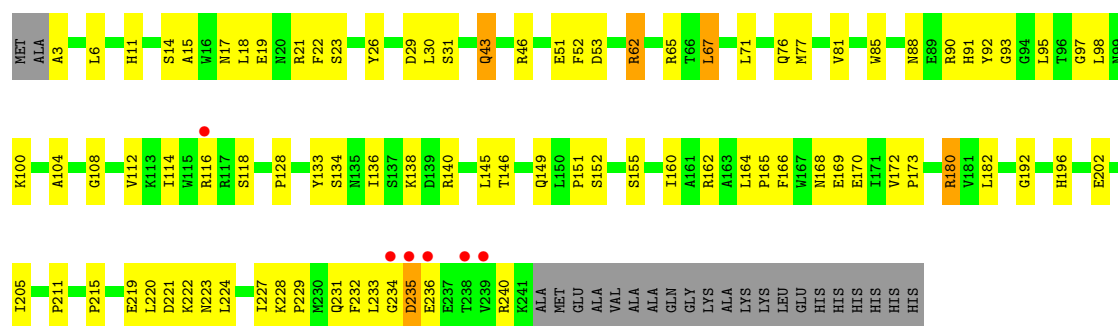


• Molecule 1: Phosphoglycerate mutase 1

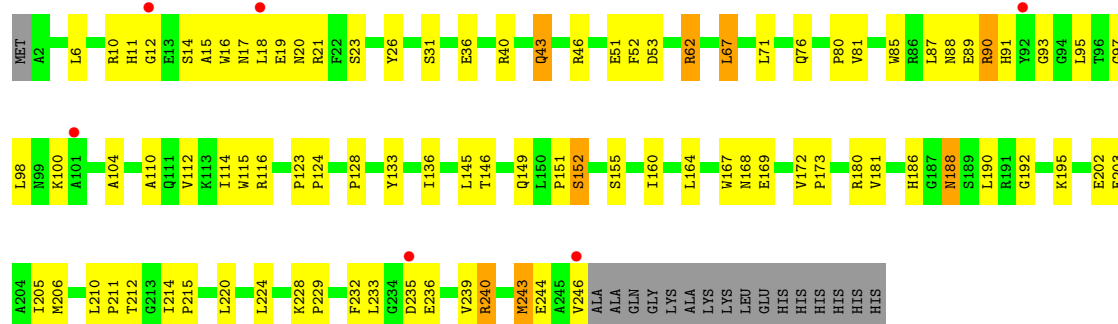




• Molecule 1: Phosphoglycerate mutase 1



• Molecule 1: Phosphoglycerate mutase 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	130.53Å 75.93Å 186.98Å 90.00° 94.42° 90.00°	Depositor
Resolution (Å)	29.91 – 2.80 29.91 – 2.80	Depositor EDS
% Data completeness (in resolution range)	93.9 (29.91-2.80) 93.8 (29.91-2.80)	Depositor EDS
R_{merge}	0.01	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.61 (at 2.80Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.233 , 0.272 0.234 , 0.271	Depositor DCC
R_{free} test set	8542 reflections (10.05%)	wwPDB-VP
Wilson B-factor (Å ²)	35.5	Xtriage
Anisotropy	0.264	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 41.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	23286	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.31% 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: CIT, 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	0.48	0/1962	0.93	6/2671 (0.2%)
1	B	0.49	0/1952	0.94	4/2655 (0.2%)
1	C	0.47	0/1947	0.94	8/2652 (0.3%)
1	D	0.45	0/1963	0.95	8/2673 (0.3%)
1	E	0.46	0/1946	0.92	8/2650 (0.3%)
1	F	0.45	0/1939	0.93	6/2641 (0.2%)
1	G	0.43	0/1936	0.93	8/2641 (0.3%)
1	H	0.44	0/1924	0.95	7/2621 (0.3%)
1	I	0.44	0/1953	0.93	8/2657 (0.3%)
1	J	0.44	0/1967	0.95	8/2676 (0.3%)
1	K	0.47	0/1919	0.92	5/2612 (0.2%)
1	L	0.48	0/1970	0.95	8/2679 (0.3%)
All	All	0.46	0/23378	0.94	84/31828 (0.3%)

There are no bond length outliers.

The worst 5 of 84 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	241	LYS	N-CA-C	-10.26	100.09	111.28
1	H	117	ARG	N-CA-C	7.81	123.31	112.45
1	E	19	GLU	N-CA-C	-7.58	103.62	113.17
1	G	19	GLU	N-CA-C	-7.46	103.91	113.16
1	A	19	GLU	N-CA-C	-7.30	103.97	113.17

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	1913	0	1848	71	0
1	B	1903	0	1845	69	0
1	C	1898	0	1827	68	0
1	D	1914	0	1843	73	0
1	E	1897	0	1822	63	0
1	F	1890	0	1819	79	0
1	G	1887	0	1786	89	0
1	H	1875	0	1800	87	0
1	I	1904	0	1844	100	0
1	J	1918	0	1858	84	0
1	K	1870	0	1806	76	0
1	L	1921	0	1865	76	0
2	A	1	0	0	0	0
2	C	1	0	0	0	0
2	F	1	0	0	0	0
2	H	1	0	0	0	0
2	J	1	0	0	0	0
2	K	1	0	0	0	0
3	A	13	0	5	1	0
3	B	13	0	5	2	0
3	C	13	0	5	1	0
3	D	13	0	5	2	0
3	E	13	0	5	0	0
3	F	13	0	5	1	0
3	G	13	0	5	1	0
3	H	13	0	5	2	0
3	I	13	0	5	1	0
3	J	13	0	5	2	0
3	K	13	0	5	3	0
3	L	13	0	5	3	0
4	A	43	0	0	2	0
4	B	45	0	0	1	0
4	C	21	0	0	1	0
4	D	33	0	0	3	0
4	E	17	0	0	3	0
4	F	27	0	0	5	0
4	G	12	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	H	18	0	0	2	0
4	I	19	0	0	2	0
4	J	15	0	0	1	0
4	K	35	0	0	9	0
4	L	49	0	0	3	0
All	All	23286	0	22023	898	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 20.

The worst 5 of 898 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:195:LYS:HB2	1:A:205:ILE:HD12	1.22	1.11
1:H:195:LYS:HB2	1:H:205:ILE:HD12	1.34	1.09
1:F:206:MET:HE2	1:F:206:MET:HA	1.37	1.02
1:A:91:HIS:HD2	1:A:93:GLY:H	1.12	0.98
1:I:43:GLN:HE21	1:I:46:ARG:HH22	1.10	0.96

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	243/262 (93%)	223 (92%)	19 (8%)	1 (0%)	30	60
1	B	240/262 (92%)	220 (92%)	20 (8%)	0	100	100
1	C	241/262 (92%)	221 (92%)	20 (8%)	0	100	100
1	D	243/262 (93%)	226 (93%)	15 (6%)	2 (1%)	16	44
1	E	241/262 (92%)	224 (93%)	17 (7%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	240/262 (92%)	221 (92%)	19 (8%)	0	100	100
1	G	243/262 (93%)	221 (91%)	18 (7%)	4 (2%)	7	27
1	H	239/262 (91%)	221 (92%)	16 (7%)	2 (1%)	16	44
1	I	240/262 (92%)	222 (92%)	17 (7%)	1 (0%)	30	60
1	J	242/262 (92%)	227 (94%)	15 (6%)	0	100	100
1	K	237/262 (90%)	225 (95%)	10 (4%)	2 (1%)	16	44
1	L	243/262 (93%)	224 (92%)	18 (7%)	1 (0%)	30	60
All	All	2892/3144 (92%)	2675 (92%)	204 (7%)	13 (0%)	30	60

5 of 13 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	I	235	ASP
1	D	243	MET
1	G	110	ALA
1	G	146	THR
1	K	235	ASP

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	194/220 (88%)	185 (95%)	9 (5%)	24	58
1	B	195/220 (89%)	186 (95%)	9 (5%)	24	58
1	C	193/220 (88%)	185 (96%)	8 (4%)	27	62
1	D	194/220 (88%)	186 (96%)	8 (4%)	27	62
1	E	192/220 (87%)	186 (97%)	6 (3%)	35	70
1	F	192/220 (87%)	185 (96%)	7 (4%)	31	66
1	G	188/220 (86%)	178 (95%)	10 (5%)	20	52
1	H	190/220 (86%)	182 (96%)	8 (4%)	26	61
1	I	196/220 (89%)	187 (95%)	9 (5%)	24	58

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	J	196/220 (89%)	186 (95%)	10 (5%)	21	54
1	K	189/220 (86%)	181 (96%)	8 (4%)	26	61
1	L	195/220 (89%)	188 (96%)	7 (4%)	31	66
All	All	2314/2640 (88%)	2215 (96%)	99 (4%)	26	60

5 of 99 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	6	LEU
1	I	112	VAL
1	H	62	ARG
1	I	6	LEU
1	J	6	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 83 such sidechains are listed below:

Mol	Chain	Res	Type
1	I	17	ASN
1	K	43	GLN
1	I	76	GLN
1	J	20	ASN
1	K	168	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry

Of 18 ligands modelled in this entry, 6 are monoatomic - leaving 12 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	CIT	A	601	-	12,12,12	2.35	4 (33%)	17,17,17	3.02	8 (47%)
3	CIT	G	607	-	12,12,12	2.56	5 (41%)	17,17,17	2.86	7 (41%)
3	CIT	H	608	-	12,12,12	2.22	4 (33%)	17,17,17	2.95	9 (52%)
3	CIT	J	610	-	12,12,12	2.43	4 (33%)	17,17,17	2.60	7 (41%)
3	CIT	K	611	-	12,12,12	2.11	3 (25%)	17,17,17	3.02	10 (58%)
3	CIT	D	604	-	12,12,12	2.44	4 (33%)	17,17,17	2.52	9 (52%)
3	CIT	B	602	-	12,12,12	2.08	4 (33%)	17,17,17	3.09	9 (52%)
3	CIT	I	609	-	12,12,12	2.10	2 (16%)	17,17,17	3.05	10 (58%)
3	CIT	E	605	-	12,12,12	2.46	5 (41%)	17,17,17	2.80	9 (52%)
3	CIT	F	606	-	12,12,12	1.97	5 (41%)	17,17,17	3.02	10 (58%)
3	CIT	L	612	-	12,12,12	2.47	4 (33%)	17,17,17	2.85	8 (47%)
3	CIT	C	603	-	12,12,12	2.09	3 (25%)	17,17,17	3.12	10 (58%)

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	CIT	A	601	-	-	4/16/16/16	-
3	CIT	G	607	-	-	5/16/16/16	-
3	CIT	H	608	-	-	9/16/16/16	-
3	CIT	J	610	-	-	9/16/16/16	-
3	CIT	K	611	-	-	5/16/16/16	-
3	CIT	D	604	-	-	5/16/16/16	-
3	CIT	B	602	-	-	9/16/16/16	-
3	CIT	I	609	-	-	7/16/16/16	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	CIT	E	605	-	-	6/16/16/16	-
3	CIT	F	606	-	-	6/16/16/16	-
3	CIT	L	612	-	-	4/16/16/16	-
3	CIT	C	603	-	-	9/16/16/16	-

The worst 5 of 47 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	L	612	CIT	C3-C6	6.02	1.59	1.53
3	G	607	CIT	C3-C6	5.85	1.59	1.53
3	H	608	CIT	C3-C6	5.66	1.59	1.53
3	J	610	CIT	C3-C6	5.59	1.59	1.53
3	D	604	CIT	C3-C6	5.55	1.59	1.53

The worst 5 of 106 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	601	CIT	O7-C3-C6	-7.17	98.79	108.96
3	B	602	CIT	C2-C3-C6	6.63	124.70	110.03
3	L	612	CIT	O7-C3-C6	-6.60	99.60	108.96
3	H	608	CIT	C2-C3-C6	6.47	124.34	110.03
3	C	603	CIT	C2-C3-C6	6.17	123.68	110.03

There are no chirality outliers.

5 of 78 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	602	CIT	O7-C3-C4-C5
3	B	602	CIT	C2-C3-C6-O5
3	B	602	CIT	C2-C3-C6-O6
3	B	602	CIT	O7-C3-C6-O5
3	B	602	CIT	O7-C3-C6-O6

There are no ring outliers.

11 monomers are involved in 19 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	601	CIT	1	0
3	G	607	CIT	1	0
3	H	608	CIT	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	J	610	CIT	2	0
3	K	611	CIT	3	0
3	D	604	CIT	2	0
3	B	602	CIT	2	0
3	I	609	CIT	1	0
3	F	606	CIT	1	0
3	L	612	CIT	3	0
3	C	603	CIT	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](#)

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	245/262 (93%)	0.09	5 (2%) 65 56	19, 33, 55, 81	0
1	B	242/262 (92%)	0.02	2 (0%) 82 75	18, 33, 49, 78	0
1	C	243/262 (92%)	0.05	5 (2%) 63 54	18, 34, 51, 84	0
1	D	245/262 (93%)	0.06	2 (0%) 82 75	18, 36, 54, 74	0
1	E	243/262 (92%)	0.16	5 (2%) 63 54	21, 36, 56, 88	0
1	F	242/262 (92%)	0.11	3 (1%) 76 68	20, 36, 56, 95	0
1	G	245/262 (93%)	0.37	11 (4%) 38 30	21, 38, 61, 85	0
1	H	241/262 (91%)	0.28	3 (1%) 76 68	21, 37, 53, 90	0
1	I	242/262 (92%)	0.29	4 (1%) 69 60	20, 37, 60, 91	0
1	J	244/262 (93%)	0.23	8 (3%) 49 39	20, 38, 58, 91	0
1	K	239/262 (91%)	0.09	6 (2%) 58 48	19, 35, 49, 102	0
1	L	245/262 (93%)	0.24	6 (2%) 59 49	18, 32, 52, 82	0
All	All	2916/3144 (92%)	0.17	60 (2%) 63 54	18, 35, 57, 102	0

The worst 5 of 60 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	K	235	ASP	5.4
1	A	19	GLU	4.8
1	I	245	ALA	4.5
1	K	234	GLY	4.3
1	E	19	GLU	4.0

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 oligosaccharides 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($\text{\AA}^2$)	Q<0.9
3	CIT	J	610	13/13	0.75	0.19	64,65,69,71	0
3	CIT	G	607	13/13	0.78	0.18	72,74,77,77	0
3	CIT	C	603	13/13	0.80	0.19	55,57,59,59	0
3	CIT	H	608	13/13	0.81	0.23	72,74,81,81	0
3	CIT	D	604	13/13	0.83	0.19	65,67,70,71	0
3	CIT	K	611	13/13	0.83	0.18	61,62,63,64	0
3	CIT	I	609	13/13	0.84	0.16	55,56,58,58	0
3	CIT	F	606	13/13	0.85	0.20	73,76,81,82	0
3	CIT	B	602	13/13	0.86	0.20	69,70,73,74	0
3	CIT	E	605	13/13	0.86	0.21	69,71,72,72	0
3	CIT	L	612	13/13	0.86	0.15	49,52,55,56	0
3	CIT	A	601	13/13	0.89	0.16	48,50,55,58	0
2	CL	H	263	1/1	0.93	0.06	25,25,25,25	0
2	CL	F	263	1/1	0.95	0.06	31,31,31,31	0
2	CL	K	263	1/1	0.96	0.05	21,21,21,21	0
2	CL	J	263	1/1	0.97	0.04	32,32,32,32	0
2	CL	A	263	1/1	0.97	0.03	22,22,22,22	0
2	CL	C	263	1/1	0.98	0.04	23,23,23,23	0

6.5 Other polymers [i](#)

There are no such residues in this entry.