



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 02:31 AM UTC

PDB ID : 5BT8 / pdb_00005bt8
Title : X-ray Crystal Structure of phosphoglycerate kinase from *Acinetobacter baumannii*
Authors : Seattle Structural Genomics Center for Infectious Disease (SSGCID)
Deposited on : 2015-06-02
Resolution : 2.30 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

<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
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

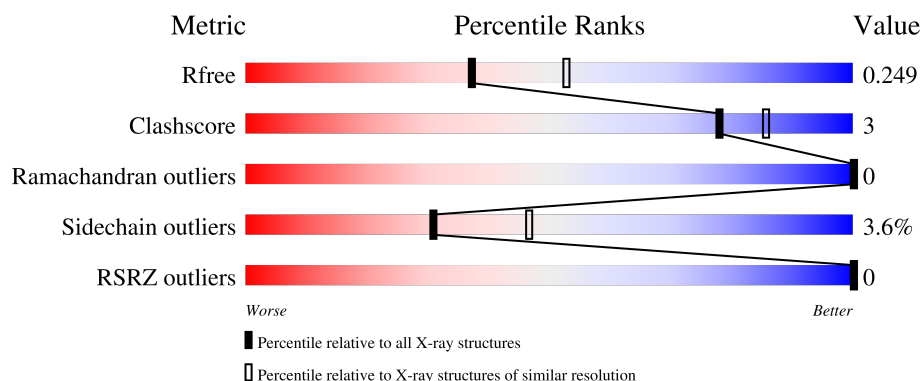
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	403	
1	B	403	
1	C	403	
1	D	403	
1	E	403	

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Mol	Chain	Length	Quality of chain
1	F	403	 91% 7% •

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 17449 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 kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	396	Total	C	N	O	S	0	1	0
			2845	1809	478	550	8			
1	B	397	Total	C	N	O	S	0	0	0
			2862	1815	482	556	9			
1	C	396	Total	C	N	O	S	0	1	0
			2843	1809	475	550	9			
1	D	395	Total	C	N	O	S	0	0	0
			2835	1803	472	551	9			
1	E	395	Total	C	N	O	S	0	0	0
			2822	1796	468	549	9			
1	F	397	Total	C	N	O	S	0	0	0
			2832	1805	476	542	9			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	MET	-	initiating methionine	UNP A0A0A2T205
A	-6	ALA	-	expression tag	UNP A0A0A2T205
A	-5	HIS	-	expression tag	UNP A0A0A2T205
A	-4	HIS	-	expression tag	UNP A0A0A2T205
A	-3	HIS	-	expression tag	UNP A0A0A2T205
A	-2	HIS	-	expression tag	UNP A0A0A2T205
A	-1	HIS	-	expression tag	UNP A0A0A2T205
A	0	HIS	-	expression tag	UNP A0A0A2T205
B	-7	MET	-	initiating methionine	UNP A0A0A2T205
B	-6	ALA	-	expression tag	UNP A0A0A2T205
B	-5	HIS	-	expression tag	UNP A0A0A2T205
B	-4	HIS	-	expression tag	UNP A0A0A2T205
B	-3	HIS	-	expression tag	UNP A0A0A2T205
B	-2	HIS	-	expression tag	UNP A0A0A2T205
B	-1	HIS	-	expression tag	UNP A0A0A2T205
B	0	HIS	-	expression tag	UNP A0A0A2T205
C	-7	MET	-	initiating methionine	UNP A0A0A2T205

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-6	ALA	-	expression tag	UNP A0A0A2T205
C	-5	HIS	-	expression tag	UNP A0A0A2T205
C	-4	HIS	-	expression tag	UNP A0A0A2T205
C	-3	HIS	-	expression tag	UNP A0A0A2T205
C	-2	HIS	-	expression tag	UNP A0A0A2T205
C	-1	HIS	-	expression tag	UNP A0A0A2T205
C	0	HIS	-	expression tag	UNP A0A0A2T205
D	-7	MET	-	initiating methionine	UNP A0A0A2T205
D	-6	ALA	-	expression tag	UNP A0A0A2T205
D	-5	HIS	-	expression tag	UNP A0A0A2T205
D	-4	HIS	-	expression tag	UNP A0A0A2T205
D	-3	HIS	-	expression tag	UNP A0A0A2T205
D	-2	HIS	-	expression tag	UNP A0A0A2T205
D	-1	HIS	-	expression tag	UNP A0A0A2T205
D	0	HIS	-	expression tag	UNP A0A0A2T205
E	-7	MET	-	initiating methionine	UNP A0A0A2T205
E	-6	ALA	-	expression tag	UNP A0A0A2T205
E	-5	HIS	-	expression tag	UNP A0A0A2T205
E	-4	HIS	-	expression tag	UNP A0A0A2T205
E	-3	HIS	-	expression tag	UNP A0A0A2T205
E	-2	HIS	-	expression tag	UNP A0A0A2T205
E	-1	HIS	-	expression tag	UNP A0A0A2T205
E	0	HIS	-	expression tag	UNP A0A0A2T205
F	-7	MET	-	initiating methionine	UNP A0A0A2T205
F	-6	ALA	-	expression tag	UNP A0A0A2T205
F	-5	HIS	-	expression tag	UNP A0A0A2T205
F	-4	HIS	-	expression tag	UNP A0A0A2T205
F	-3	HIS	-	expression tag	UNP A0A0A2T205
F	-2	HIS	-	expression tag	UNP A0A0A2T205
F	-1	HIS	-	expression tag	UNP A0A0A2T205
F	0	HIS	-	expression tag	UNP A0A0A2T205

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	69	Total O 69 69	0	0
2	B	74	Total O 74 74	0	0
2	C	66	Total O 66 66	0	0
2	D	67	Total O 67 67	0	0

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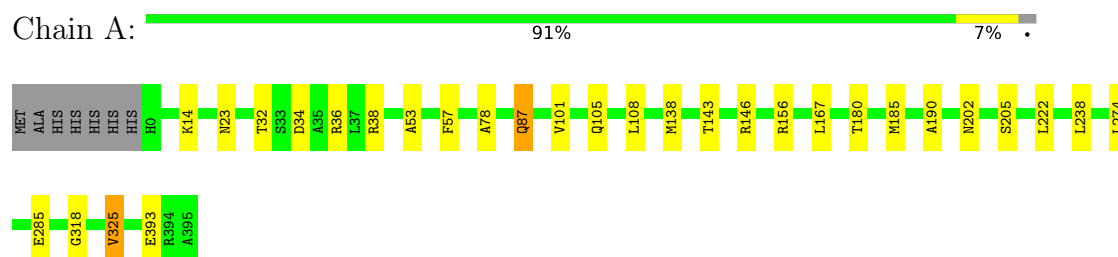
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	E	89	Total	O	0	0
			89	89		
2	F	45	Total	O	0	0
			45	45		

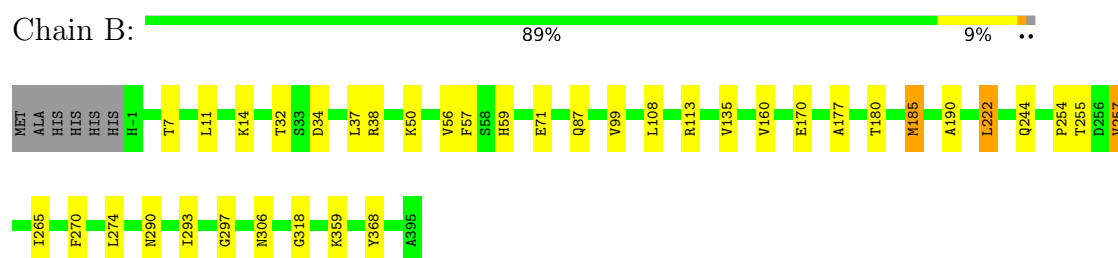
3 Residue-property plots [i](#)

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.

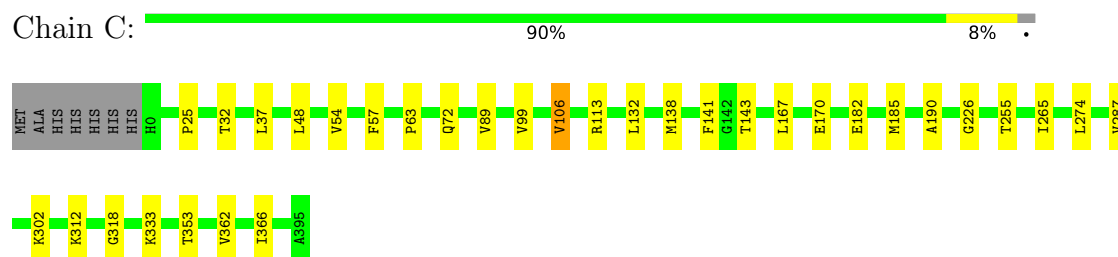
- Molecule 1: Phosphoglycerate kinase



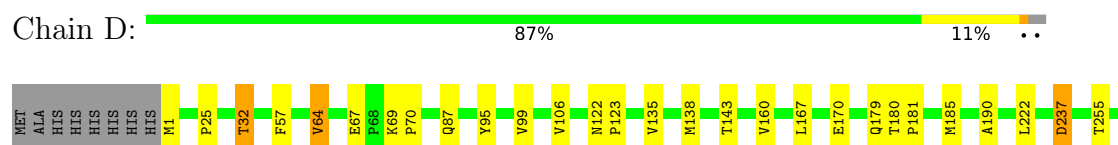
- Molecule 1: Phosphoglycerate kinase

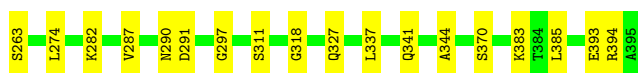


- Molecule 1: Phosphoglycerate kinase



- Molecule 1: Phosphoglycerate kinase





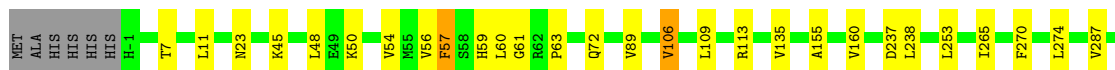
- Molecule 1: Phosphoglycerate kinase

Chain E: 88% 10% .



- Molecule 1: Phosphoglycerate kinase

Chain F: 91% 7% .



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	70.44Å 200.29Å 87.30Å 90.00° 91.37° 90.00°	Depositor
Resolution (Å)	48.45 – 2.30 48.45 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.45-2.30) 99.9 (48.45-2.30)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.49 (at 2.29Å)	Xtriage
Refinement program	PHENIX (dev_2050)	Depositor
R, R_{free}	0.215 , 0.248 0.217 , 0.249	Depositor DCC
R_{free} test set	2003 reflections (1.87%)	wwPDB-VP
Wilson B-factor (Å ²)	33.2	Xtriage
Anisotropy	0.565	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 14.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.095 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	17449	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.85% 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

5.1 Standard geometry

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	0.33	0/2891	0.75	0/3945
1	B	0.34	0/2907	0.74	2/3965 (0.1%)
1	C	0.34	0/2890	0.73	1/3943 (0.0%)
1	D	0.34	0/2878	0.76	2/3927 (0.1%)
1	E	0.34	0/2865	0.74	2/3911 (0.1%)
1	F	0.34	0/2877	0.74	2/3928 (0.1%)
All	All	0.34	0/17308	0.74	9/23619 (0.0%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	297	GLY	CA-C-N	6.03	125.24	118.97
1	B	297	GLY	C-N-CA	6.03	125.24	118.97
1	C	255	THR	N-CA-C	-5.63	106.58	113.50
1	F	297	GLY	CA-C-N	5.52	124.98	119.24
1	F	297	GLY	C-N-CA	5.52	124.98	119.24
1	E	297	GLY	CA-C-N	5.19	125.24	119.32
1	E	297	GLY	C-N-CA	5.19	125.24	119.32
1	D	297	GLY	CA-C-N	5.18	125.23	119.32
1	D	297	GLY	C-N-CA	5.18	125.23	119.32

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	2845	0	2830	11	0
1	B	2862	0	2837	14	0
1	C	2843	0	2820	14	0
1	D	2835	0	2812	22	0
1	E	2822	0	2792	17	0
1	F	2832	0	2800	19	0
2	A	69	0	0	1	0
2	B	74	0	0	0	0
2	C	66	0	0	0	0
2	D	67	0	0	2	0
2	E	89	0	0	0	0
2	F	45	0	0	1	0
All	All	17449	0	16891	95	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 (95) 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:257:VAL:HG11	1:B:293:ILE:HG12	1.71	0.71
1:E:60:LEU:HD12	1:E:74:LEU:HD23	1.73	0.69
1:E:257:VAL:HG11	1:E:293:ILE:HG12	1.77	0.66
1:D:1:MET:HE2	1:D:179:GLN:HE22	1.62	0.65
1:D:274:LEU:O	1:D:327:GLN:NE2	2.29	0.65
1:F:265:ILE:HD12	1:F:274:LEU:HD12	1.77	0.65
1:D:237:ASP:OD1	1:D:237:ASP:N	2.25	0.65
1:F:54:VAL:HB	1:F:106:VAL:HG22	1.80	0.64
1:A:34:ASP:OD2	1:A:38:ARG:NH1	2.30	0.63
1:F:23:ASN:HA	1:F:60:LEU:HD12	1.82	0.61
1:B:265:ILE:HD12	1:B:274:LEU:HD12	1.83	0.61
1:F:135:VAL:HG22	1:F:160:VAL:HB	1.83	0.61
1:B:34:ASP:OD1	1:B:38:ARG:NH1	2.34	0.60
1:D:64:VAL:HG13	1:D:67:GLU:HB2	1.82	0.59
1:F:89:VAL:HG13	1:F:106:VAL:HG12	1.84	0.59
1:E:282:LYS:NZ	1:E:291:ASP:OD2	2.29	0.58
1:D:393:GLU:OE1	2:D:401:HOH:O	2.17	0.58
1:D:337:LEU:O	1:D:341:GLN:HG2	2.04	0.58
1:D:394:ARG:NH2	2:D:408:HOH:O	2.38	0.56
1:D:282:LYS:NZ	1:D:291:ASP:OD2	2.30	0.56
1:D:95:TYR:HB2	1:D:99:VAL:HG22	1.89	0.54
1:C:48:LEU:HD21	1:C:106:VAL:HG23	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:54:VAL:HB	1:C:106:VAL:HG22	1.92	0.52
1:D:138:MET:HG3	1:D:167:LEU:HD22	1.93	0.51
1:C:190:ALA:HB3	1:C:318:GLY:HA3	1.93	0.51
1:C:182:GLU:HG2	1:C:312:LYS:HB3	1.93	0.51
1:A:274:LEU:HD23	1:A:325:VAL:HG11	1.92	0.51
1:B:190:ALA:HB3	1:B:318:GLY:HA3	1.93	0.51
1:C:89:VAL:HG13	1:C:106:VAL:HG12	1.92	0.51
1:E:219:ASN:HA	1:E:222:LEU:HD12	1.93	0.51
1:A:87:GLN:HE22	1:A:105:GLN:HA	1.76	0.50
1:A:14:LYS:HE3	2:A:422:HOH:O	2.12	0.50
1:B:222:LEU:HD11	1:B:254:PRO:HD2	1.93	0.49
1:E:95:TYR:HB2	1:E:99:VAL:HG22	1.94	0.49
1:E:274:LEU:HB3	1:E:325:VAL:HG21	1.94	0.48
1:C:265:ILE:HD12	1:C:274:LEU:HD12	1.94	0.48
1:E:190:ALA:HB3	1:E:318:GLY:HA3	1.95	0.48
1:A:156:ARG:NH1	1:A:393:GLU:OE1	2.46	0.48
1:B:59:HIS:HB2	1:B:113:ARG:HG3	1.95	0.47
1:D:263:SER:OG	1:D:290:ASN:ND2	2.47	0.47
1:B:270:PHE:O	1:B:274:LEU:HB2	2.15	0.47
1:C:99:VAL:HG12	1:C:132:LEU:HD22	1.95	0.47
1:A:202:ASN:O	1:A:205:SER:OG	2.26	0.47
1:F:48:LEU:HD21	1:F:106:VAL:HG23	1.97	0.47
1:A:23:ASN:O	1:A:36:ARG:NH1	2.43	0.47
1:F:364:ASP:OD1	1:F:364:ASP:N	2.47	0.47
1:E:25:PRO:HB2	1:E:32:THR:OG1	2.16	0.46
1:B:290:ASN:OD1	1:E:8:ASP:HB3	2.16	0.46
1:C:226:GLY:HA2	1:D:180:THR:HG22	1.98	0.45
1:F:270:PHE:O	1:F:274:LEU:HB2	2.16	0.45
1:C:25:PRO:HB2	1:C:32:THR:OG1	2.17	0.45
1:D:190:ALA:HB3	1:D:318:GLY:HA3	1.98	0.45
1:C:138:MET:HG3	1:C:167:LEU:HD22	1.99	0.45
1:F:57:PHE:H	1:F:57:PHE:HD1	1.65	0.44
1:C:362:VAL:O	1:C:366:ILE:HG12	2.18	0.44
1:B:11:LEU:HA	1:B:14:LYS:HG3	1.98	0.44
1:F:237:ASP:OD1	1:F:237:ASP:N	2.49	0.44
1:D:181:PRO:HG3	1:D:185:MET:HE2	1.99	0.44
1:E:57:PHE:HB3	1:E:109:LEU:HB2	2.00	0.44
1:E:270:PHE:O	1:E:274:LEU:HB2	2.17	0.44
1:F:61:GLY:C	1:F:63:PRO:HD3	2.43	0.44
1:E:56:VAL:O	1:E:108:LEU:HD12	2.17	0.44
1:D:383:LYS:HB3	1:D:383:LYS:HE2	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:7:THR:HA	1:F:50:LYS:HE3	2.01	0.43
1:A:53:ALA:HB1	1:A:101:VAL:HG22	2.01	0.43
1:A:78:ALA:HB2	1:A:108:LEU:HD23	2.01	0.43
1:E:170:GLU:OE1	1:E:370:SER:OG	2.36	0.43
1:F:45:LYS:NZ	2:F:407:HOH:O	2.52	0.43
1:D:135:VAL:HG22	1:D:160:VAL:HB	2.01	0.42
1:B:7:THR:HA	1:B:50:LYS:HE3	2.02	0.42
1:B:170:GLU:OE1	1:B:368:TYR:OH	2.31	0.42
1:F:155:ALA:O	1:F:394:ARG:NH1	2.52	0.42
1:F:57:PHE:HB3	1:F:109:LEU:HB2	2.01	0.42
1:D:222:LEU:HD23	1:D:222:LEU:HA	1.82	0.42
1:E:239:VAL:O	1:E:243:LYS:HG3	2.20	0.42
1:F:72:GLN:HE21	1:F:72:GLN:HB2	1.59	0.42
1:A:190:ALA:HB3	1:A:318:GLY:HA3	2.01	0.42
1:F:253:LEU:HD23	1:F:253:LEU:HA	1.89	0.42
1:C:318:GLY:N	1:C:353:THR:OG1	2.53	0.41
1:E:122:ASN:HA	1:E:123:PRO:HD3	1.95	0.41
1:E:292:MET:HE2	1:E:292:MET:HB3	1.95	0.41
1:B:177:ALA:O	1:B:185:MET:HE2	2.20	0.41
1:C:63:PRO:HB3	1:C:72:GLN:OE1	2.20	0.41
1:D:25:PRO:HB2	1:D:32:THR:OG1	2.21	0.41
1:D:170:GLU:OE1	1:D:370:SER:OG	2.31	0.41
1:F:11:LEU:HD23	1:F:50:LYS:HD3	2.02	0.41
1:A:138:MET:HG3	1:A:167:LEU:HD22	2.03	0.41
1:D:122:ASN:HA	1:D:123:PRO:HD3	1.96	0.41
1:B:135:VAL:HG22	1:B:160:VAL:HB	2.03	0.41
1:C:141:PHE:CG	1:C:170:GLU:HG2	2.56	0.41
1:F:59:HIS:HB2	1:F:113:ARG:HG2	2.03	0.41
1:B:56:VAL:O	1:B:108:LEU:HD12	2.21	0.41
1:D:311:SER:O	1:D:344:ALA:HB2	2.22	0.40
1:D:69:LYS:HA	1:D:70:PRO:HD3	1.90	0.40
1:E:99:VAL:HG23	1:E:132:LEU:HD22	2.03	0.40

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	395/403 (98%)	390 (99%)	5 (1%)	0	100	100
1	B	395/403 (98%)	387 (98%)	8 (2%)	0	100	100
1	C	395/403 (98%)	386 (98%)	9 (2%)	0	100	100
1	D	393/403 (98%)	387 (98%)	6 (2%)	0	100	100
1	E	393/403 (98%)	385 (98%)	8 (2%)	0	100	100
1	F	395/403 (98%)	387 (98%)	8 (2%)	0	100	100
All	All	2366/2418 (98%)	2322 (98%)	44 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	282/307 (92%)	271 (96%)	11 (4%)	28	43
1	B	286/307 (93%)	272 (95%)	14 (5%)	22	34
1	C	282/307 (92%)	273 (97%)	9 (3%)	34	51
1	D	281/307 (92%)	271 (96%)	10 (4%)	31	47
1	E	280/307 (91%)	268 (96%)	12 (4%)	26	39
1	F	277/307 (90%)	272 (98%)	5 (2%)	51	70
All	All	1688/1842 (92%)	1627 (96%)	61 (4%)	31	47

All (61) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	THR
1	A	57	PHE
1	A	87	GLN

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Mol	Chain	Res	Type
1	A	143	THR
1	A	146	ARG
1	A	180	THR
1	A	185	MET
1	A	222	LEU
1	A	238	LEU
1	A	285	GLU
1	A	325	VAL
1	B	32	THR
1	B	37	LEU
1	B	57	PHE
1	B	71	GLU
1	B	87	GLN
1	B	99	VAL
1	B	180	THR
1	B	185	MET
1	B	222	LEU
1	B	244	GLN
1	B	255	THR
1	B	257	VAL
1	B	306	ASN
1	B	359	LYS
1	C	37	LEU
1	C	57	PHE
1	C	106	VAL
1	C	113	ARG
1	C	143	THR
1	C	185	MET
1	C	287	VAL
1	C	302	LYS
1	C	333	LYS
1	D	32	THR
1	D	57	PHE
1	D	64	VAL
1	D	87	GLN
1	D	106	VAL
1	D	143	THR
1	D	237	ASP
1	D	255	THR
1	D	287	VAL
1	D	385	LEU
1	E	1	MET

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Mol	Chain	Res	Type
1	E	22	LEU
1	E	57	PHE
1	E	62	ARG
1	E	87	GLN
1	E	88	GLU
1	E	185	MET
1	E	196	THR
1	E	255	THR
1	E	257	VAL
1	E	306	ASN
1	E	364	ASP
1	F	56	VAL
1	F	57	PHE
1	F	106	VAL
1	F	238	LEU
1	F	287	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	4	GLN
1	A	72	GLN
1	A	87	GLN
1	A	210	GLN
1	B	87	GLN
1	B	122	ASN
1	B	266	ASN
1	C	4	GLN
1	C	10	ASN
1	C	115	ASN
1	D	179	GLN
1	D	210	GLN
1	D	278	GLN
1	D	290	ASN
1	D	317	ASN
1	E	4	GLN
1	E	10	ASN
1	E	219	ASN
1	E	327	GLN
1	F	72	GLN
1	F	228	ASN
1	F	327	GLN

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Mol	Chain	Res	Type
1	F	341	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	396/403 (98%)	-1.12	0 100 100	19, 42, 59, 76	1 (0%)
1	B	397/403 (98%)	-1.19	0 100 100	23, 38, 55, 65	0
1	C	396/403 (98%)	-1.14	0 100 100	21, 39, 58, 89	1 (0%)
1	D	395/403 (98%)	-1.08	0 100 100	24, 38, 66, 76	0
1	E	395/403 (98%)	-1.16	0 100 100	21, 36, 57, 67	0
1	F	397/403 (98%)	-1.11	0 100 100	24, 42, 60, 74	0
All	All	2376/2418 (98%)	-1.13	0 100 100	19, 40, 60, 89	2 (0%)

There are no RSRZ outliers to report.

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.