



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 03:22 PM GMT

PDB ID : 2Q5R  
Title : Structure of apo Staphylococcus aureus D-tagatose-6-phosphatekinase  
Authors : McGrath, T.E.; Soloveychik, M.; Romanov, V.; Thambipillai, D.; Dharamsi, A.; Virag, C.; Domagala, M.; Pai, E.F.; Edwards, A.M.; Battaile, K.; Chirgadze, N.Y.  
Deposited on : 2007-06-01  
Resolution : 2.30 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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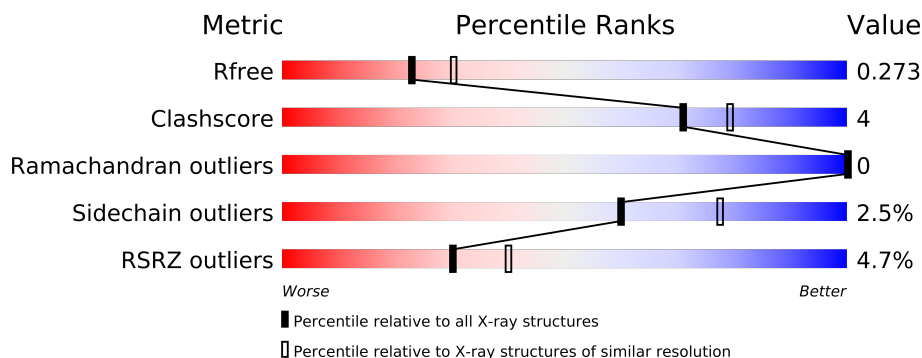
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

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))
$R_{free}$	66092	2929 (2.30-2.30)
Clashscore	79885	3679 (2.30-2.30)
Ramachandran outliers	78287	3642 (2.30-2.30)
Sidechain outliers	78261	3641 (2.30-2.30)
RSRZ outliers	66119	2930 (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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	330	
1	B	330	
1	C	330	
1	D	330	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10117 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Tagatose-6-phosphate kinase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	318	Total	C	N	O	S	Se	0	3	0
			2459	1560	416	478	3	2			
1	B	312	Total	C	N	O	S	Se	0	2	0
			2413	1531	406	471	3	2			
1	C	303	Total	C	N	O	S	Se	0	3	0
			2350	1494	395	456	3	2			
1	D	310	Total	C	N	O	S	Se	0	2	0
			2396	1519	404	467	3	3			

There are 88 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MSE	-	CLONING ARTIFACT	UNP Q5HE12
A	-18	GLY	-	CLONING ARTIFACT	UNP Q5HE12
A	-17	SER	-	CLONING ARTIFACT	UNP Q5HE12
A	-16	SER	-	CLONING ARTIFACT	UNP Q5HE12
A	-15	HIS	-	CLONING ARTIFACT	UNP Q5HE12
A	-14	HIS	-	CLONING ARTIFACT	UNP Q5HE12
A	-13	HIS	-	CLONING ARTIFACT	UNP Q5HE12
A	-12	HIS	-	CLONING ARTIFACT	UNP Q5HE12
A	-11	HIS	-	CLONING ARTIFACT	UNP Q5HE12
A	-10	HIS	-	CLONING ARTIFACT	UNP Q5HE12
A	-9	SER	-	CLONING ARTIFACT	UNP Q5HE12
A	-8	SER	-	CLONING ARTIFACT	UNP Q5HE12
A	-7	GLY	-	CLONING ARTIFACT	UNP Q5HE12
A	-6	LEU	-	CLONING ARTIFACT	UNP Q5HE12
A	-5	VAL	-	CLONING ARTIFACT	UNP Q5HE12
A	-4	PRO	-	CLONING ARTIFACT	UNP Q5HE12
A	-3	ARG	-	CLONING ARTIFACT	UNP Q5HE12
A	-2	GLY	-	CLONING ARTIFACT	UNP Q5HE12
A	-1	SER	-	CLONING ARTIFACT	UNP Q5HE12
A	0	HIS	-	CLONING ARTIFACT	UNP Q5HE12
A	1	MSE	MET	MODIFIED RESIDUE	UNP Q5HE12

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Chain	Residue	Modelled	Actual	Comment	Reference
A	282	MSE	MET	MODIFIED RESIDUE	UNP Q5HE12
B	-19	MSE	-	CLONING ARTIFACT	UNP Q5HE12
B	-18	GLY	-	CLONING ARTIFACT	UNP Q5HE12
B	-17	SER	-	CLONING ARTIFACT	UNP Q5HE12
B	-16	SER	-	CLONING ARTIFACT	UNP Q5HE12
B	-15	HIS	-	CLONING ARTIFACT	UNP Q5HE12
B	-14	HIS	-	CLONING ARTIFACT	UNP Q5HE12
B	-13	HIS	-	CLONING ARTIFACT	UNP Q5HE12
B	-12	HIS	-	CLONING ARTIFACT	UNP Q5HE12
B	-11	HIS	-	CLONING ARTIFACT	UNP Q5HE12
B	-10	HIS	-	CLONING ARTIFACT	UNP Q5HE12
B	-9	SER	-	CLONING ARTIFACT	UNP Q5HE12
B	-8	SER	-	CLONING ARTIFACT	UNP Q5HE12
B	-7	GLY	-	CLONING ARTIFACT	UNP Q5HE12
B	-6	LEU	-	CLONING ARTIFACT	UNP Q5HE12
B	-5	VAL	-	CLONING ARTIFACT	UNP Q5HE12
B	-4	PRO	-	CLONING ARTIFACT	UNP Q5HE12
B	-3	ARG	-	CLONING ARTIFACT	UNP Q5HE12
B	-2	GLY	-	CLONING ARTIFACT	UNP Q5HE12
B	-1	SER	-	CLONING ARTIFACT	UNP Q5HE12
B	0	HIS	-	CLONING ARTIFACT	UNP Q5HE12
B	1	MSE	MET	MODIFIED RESIDUE	UNP Q5HE12
B	282	MSE	MET	MODIFIED RESIDUE	UNP Q5HE12
C	-19	MSE	-	CLONING ARTIFACT	UNP Q5HE12
C	-18	GLY	-	CLONING ARTIFACT	UNP Q5HE12
C	-17	SER	-	CLONING ARTIFACT	UNP Q5HE12
C	-16	SER	-	CLONING ARTIFACT	UNP Q5HE12
C	-15	HIS	-	CLONING ARTIFACT	UNP Q5HE12
C	-14	HIS	-	CLONING ARTIFACT	UNP Q5HE12
C	-13	HIS	-	CLONING ARTIFACT	UNP Q5HE12
C	-12	HIS	-	CLONING ARTIFACT	UNP Q5HE12
C	-11	HIS	-	CLONING ARTIFACT	UNP Q5HE12
C	-10	HIS	-	CLONING ARTIFACT	UNP Q5HE12
C	-9	SER	-	CLONING ARTIFACT	UNP Q5HE12
C	-8	SER	-	CLONING ARTIFACT	UNP Q5HE12
C	-7	GLY	-	CLONING ARTIFACT	UNP Q5HE12
C	-6	LEU	-	CLONING ARTIFACT	UNP Q5HE12
C	-5	VAL	-	CLONING ARTIFACT	UNP Q5HE12
C	-4	PRO	-	CLONING ARTIFACT	UNP Q5HE12
C	-3	ARG	-	CLONING ARTIFACT	UNP Q5HE12
C	-2	GLY	-	CLONING ARTIFACT	UNP Q5HE12
C	-1	SER	-	CLONING ARTIFACT	UNP Q5HE12

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Chain	Residue	Modelled	Actual	Comment	Reference
C	0	HIS	-	CLONING ARTIFACT	UNP Q5HE12
C	1	MSE	MET	MODIFIED RESIDUE	UNP Q5HE12
C	282	MSE	MET	MODIFIED RESIDUE	UNP Q5HE12
D	-19	MSE	-	CLONING ARTIFACT	UNP Q5HE12
D	-18	GLY	-	CLONING ARTIFACT	UNP Q5HE12
D	-17	SER	-	CLONING ARTIFACT	UNP Q5HE12
D	-16	SER	-	CLONING ARTIFACT	UNP Q5HE12
D	-15	HIS	-	CLONING ARTIFACT	UNP Q5HE12
D	-14	HIS	-	CLONING ARTIFACT	UNP Q5HE12
D	-13	HIS	-	CLONING ARTIFACT	UNP Q5HE12
D	-12	HIS	-	CLONING ARTIFACT	UNP Q5HE12
D	-11	HIS	-	CLONING ARTIFACT	UNP Q5HE12
D	-10	HIS	-	CLONING ARTIFACT	UNP Q5HE12
D	-9	SER	-	CLONING ARTIFACT	UNP Q5HE12
D	-8	SER	-	CLONING ARTIFACT	UNP Q5HE12
D	-7	GLY	-	CLONING ARTIFACT	UNP Q5HE12
D	-6	LEU	-	CLONING ARTIFACT	UNP Q5HE12
D	-5	VAL	-	CLONING ARTIFACT	UNP Q5HE12
D	-4	PRO	-	CLONING ARTIFACT	UNP Q5HE12
D	-3	ARG	-	CLONING ARTIFACT	UNP Q5HE12
D	-2	GLY	-	CLONING ARTIFACT	UNP Q5HE12
D	-1	SER	-	CLONING ARTIFACT	UNP Q5HE12
D	0	HIS	-	CLONING ARTIFACT	UNP Q5HE12
D	1	MSE	MET	MODIFIED RESIDUE	UNP Q5HE12
D	282	MSE	MET	MODIFIED RESIDUE	UNP Q5HE12

- Molecule 2 is water.

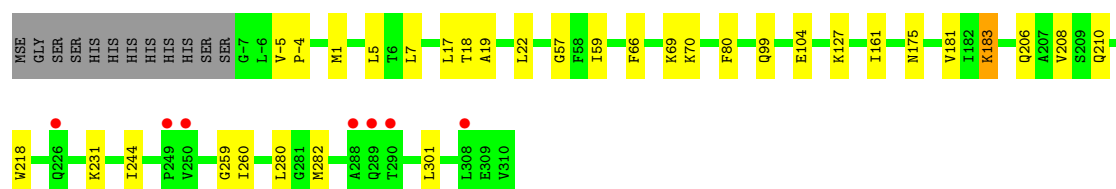
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	122	Total O 122 122	0	0
2	B	136	Total O 136 136	0	0
2	C	99	Total O 99 99	0	0
2	D	142	Total O 142 142	0	0

### 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.

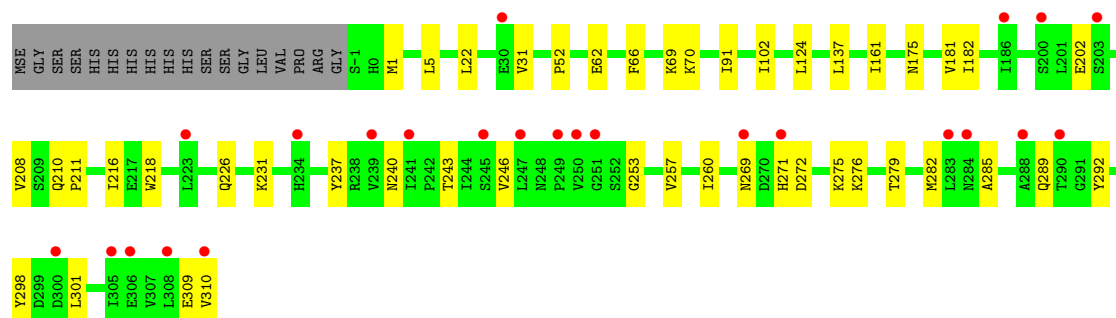
- Molecule 1: Tagatose-6-phosphate kinase

Chain A: 



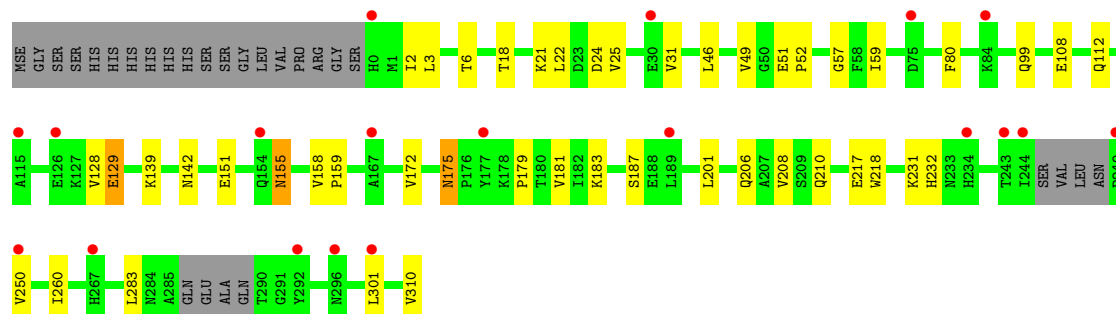
- Molecule 1: Tagatose-6-phosphate kinase

Chain B: 



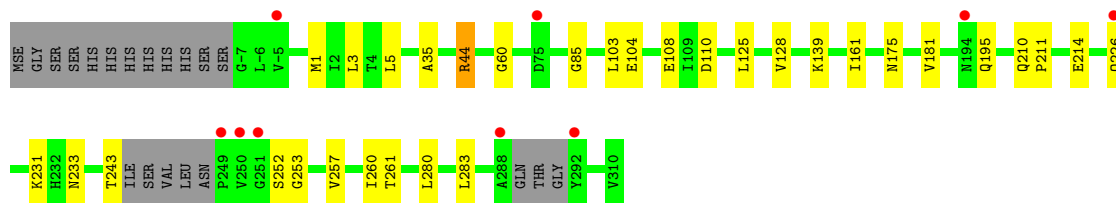
- Molecule 1: Tagatose-6-phosphate kinase

Chain C: 



- Molecule 1: Tagatose-6-phosphate kinase

Chain D: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	76.95Å 96.94Å 94.48Å 90.00° 94.42° 90.00°	Depositor
Resolution (Å)	14.99 – 2.30 14.99 – 2.30	Depositor EDS
% Data completeness (in resolution range)	90.1 (14.99-2.30) 90.1 (14.99-2.30)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.85 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.215 , 0.283 0.209 , 0.273	Depositor DCC
$R_{free}$ test set	2729 reflections (5.20%)	DCC
Wilson B-factor (Å <sup>2</sup> )	39.9	Xtriage
Anisotropy	0.053	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 52.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 55227 reflections (0.002%)	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	10117	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.63% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.



## 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	0.44	0/2508	0.58	0/3403
1	B	0.59	3/2458 (0.1%)	0.59	0/3335
1	C	0.64	4/2397 (0.2%)	0.58	0/3250
1	D	0.48	0/2440	0.59	1/3306 (0.0%)
All	All	0.54	7/9803 (0.1%)	0.59	1/13294 (0.0%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	155	ASN	CG-OD1	15.47	1.57	1.24
1	C	155	ASN	CG-ND2	14.75	1.69	1.32
1	B	202	GLU	CG-CD	6.67	1.61	1.51
1	C	151	GLU	CD-OE2	5.53	1.31	1.25
1	C	151	GLU	CD-OE1	5.46	1.31	1.25
1	B	240	ASN	C-N	5.25	1.46	1.34
1	B	310	VAL	CB-CG2	5.24	1.63	1.52

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	44	ARG	NE-CZ-NH2	-5.12	117.74	120.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the

chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2459	0	2482	25	0
1	B	2413	0	2430	24	0
1	C	2350	0	2365	28	0
1	D	2396	0	2404	15	0
2	A	122	0	0	1	0
2	B	136	0	0	0	0
2	C	99	0	0	3	0
2	D	142	0	0	2	0
All	All	10117	0	9681	86	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 4.

All (86) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:155:ASN:CG	1:C:155:ASN:ND2	1.69	1.46
1:C:129:GLU:O	1:C:159:PRO:HD2	1.70	0.91
1:A:183:LYS:O	1:A:183:LYS:HG2	1.77	0.83
1:A:244:ILE:HD13	1:A:282:MSE:HG2	1.67	0.75
1:A:244:ILE:HG21	1:A:282:MSE:HE3	1.74	0.70
1:B:181:VAL:HG11	1:B:260:ILE:HG21	1.76	0.67
1:A:59:ILE:HG12	1:A:80:PHE:HB3	1.79	0.65
1:A:66:PHE:CZ	1:A:70:LYS:HE2	2.33	0.64
1:B:66:PHE:CZ	1:B:70[B]:LYS:HE3	2.35	0.62
1:D:35:ALA:O	1:D:44:ARG:NH2	2.29	0.61
1:D:280:LEU:HD23	1:D:283:LEU:HD12	1.84	0.60
1:C:155:ASN:ND2	1:C:155:ASN:CB	2.61	0.60
1:C:31:VAL:HG21	1:D:104:GLU:HG2	1.84	0.59
1:C:6:THR:O	1:C:57:GLY:HA3	2.02	0.59
1:A:282:MSE:HE2	1:A:301:LEU:HD13	1.87	0.57
1:C:59:ILE:HG12	1:C:80:PHE:HB3	1.86	0.56
1:B:279:THR:HG23	1:B:301:LEU:HD22	1.88	0.56
1:C:181:VAL:HG22	1:C:218:TRP:HB2	1.87	0.56
1:A:17[B]:LEU:HG	1:A:19:ALA:O	2.06	0.55
1:A:22:LEU:HD11	1:B:22:LEU:HD11	1.88	0.55
1:A:69[B]:LYS:HG3	1:B:62:GLU:OE2	2.07	0.55
1:A:244:ILE:CD1	1:A:282:MSE:HG2	2.37	0.55
1:D:211:PRO:HA	1:D:214:GLU:HG3	1.89	0.54
1:A:7:LEU:HD23	1:A:57:GLY:HA2	1.90	0.53
1:A:59:ILE:HG12	1:A:80:PHE:CB	2.39	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:282:MSE:SE	1:B:301:LEU:HD21	2.60	0.52
1:D:195:GLN:NE2	2:D:437:HOH:O	2.41	0.52
1:D:1[A]:MSE:SE	2:D:403:HOH:O	2.77	0.52
1:A:181:VAL:HG11	1:A:260:ILE:HG21	1.92	0.52
1:A:244:ILE:CG2	1:A:282:MSE:HE3	2.39	0.51
1:C:51:GLU:OE1	1:C:52:PRO:HD2	2.12	0.50
1:C:108:GLU:HG3	1:C:139:LYS:HD3	1.95	0.49
1:C:129:GLU:O	1:C:159:PRO:CD	2.53	0.49
1:A:161:ILE:N	1:A:161:ILE:HD12	2.28	0.48
1:A:104:GLU:HG2	1:B:31:VAL:HG21	1.95	0.48
1:B:181:VAL:HG22	1:B:218:TRP:HB2	1.96	0.48
1:C:181:VAL:HG11	1:C:260:ILE:HG21	1.95	0.48
1:C:99:GLN:OE1	1:C:250:VAL:HG22	2.13	0.48
1:A:69[A]:LYS:HD2	1:B:62:GLU:OE2	2.13	0.48
1:C:179:PRO:HA	2:C:400:HOH:O	2.13	0.47
1:B:276:LYS:HG3	1:B:298:TYR:OH	2.15	0.47
1:B:161:ILE:HD12	1:B:161:ILE:N	2.30	0.47
1:D:108:GLU:HG3	1:D:139:LYS:HG2	1.96	0.46
1:C:250:VAL:HG21	2:C:340:HOH:O	2.15	0.46
1:B:289:GLN:HB2	1:B:292:TYR:HB3	1.97	0.46
1:A:208:VAL:HG23	1:A:231:LYS:HD2	1.96	0.46
1:C:206:GLN:O	1:C:210:GLN:HG2	2.15	0.46
1:B:246:VAL:HG12	1:B:285:ALA:HB1	1.97	0.46
1:C:283:LEU:HD12	1:C:301:LEU:HD12	1.97	0.45
1:B:237:TYR:CE1	1:B:309:GLU:HG3	2.52	0.45
1:B:271:HIS:O	1:B:275:LYS:HG3	2.17	0.45
1:C:208:VAL:HG12	1:C:231:LYS:HD3	1.98	0.44
1:D:181:VAL:HG11	1:D:260:ILE:HG21	1.98	0.44
1:C:21:LYS:HB2	1:C:21:LYS:HE3	1.81	0.44
1:C:158:VAL:HA	1:C:159:PRO:HD3	1.85	0.44
1:D:60:GLY:HA3	1:D:85:GLY:O	2.17	0.43
1:B:253:GLY:O	1:B:257:VAL:HG23	2.17	0.43
1:A:1:MSE:HE2	1:A:127:LYS:O	2.18	0.43
1:B:91[B]:ILE:HG23	1:B:102:ILE:HB	1.99	0.43
1:B:182:ILE:HG12	1:B:216:ILE:HG21	2.00	0.43
1:D:253:GLY:O	1:D:257:VAL:HG23	2.18	0.43
1:A:181:VAL:HG22	1:A:218:TRP:HB2	2.00	0.43
1:A:259:GLY:CA	1:A:280:LEU:HD12	2.49	0.43
1:D:3:LEU:HB2	1:D:128:VAL:HG11	2.00	0.43
1:B:1:MSE:HA	1:B:52:PRO:HB2	2.01	0.43
1:C:46:LEU:HA	1:C:49:VAL:HG22	2.01	0.43
1:A:99:GLN:NE2	2:A:417:HOH:O	2.52	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:210:GLN:HG3	1:B:211:PRO:HD2	2.01	0.42
1:C:217:GLU:HG2	1:C:232:HIS:CD2	2.55	0.42
1:B:69:LYS:HE2	1:B:69:LYS:HB3	1.79	0.42
1:C:172:VAL:HA	1:C:175:ASN:HD22	1.84	0.42
1:B:62:GLU:CD	1:B:62:GLU:H	2.23	0.41
1:C:112:GLN:CD	1:C:112:GLN:H	2.23	0.41
1:B:279:THR:HA	1:B:301:LEU:HD22	2.01	0.41
1:C:201:LEU:CD1	1:C:310:VAL:HG12	2.50	0.41
1:C:142:ASN:HB3	2:C:394:HOH:O	2.20	0.41
1:C:25[A]:VAL:HG13	1:D:103:LEU:HG	2.01	0.41
1:C:2:ILE:HD12	1:C:51:GLU:HB3	2.03	0.41
1:A:206:GLN:O	1:A:210:GLN:HG2	2.20	0.41
1:D:110:ASP:C	1:D:110:ASP:OD1	2.59	0.41
1:A:183:LYS:O	1:A:183:LYS:CG	2.58	0.41
1:A:-5:VAL:HA	1:A:-4:PRO:HD3	1.93	0.41
1:B:208:VAL:HG13	1:B:231:LYS:HD2	2.02	0.41
1:D:161:ILE:HD13	1:D:261:THR:HG23	2.03	0.41
1:D:231:LYS:HE2	1:D:233:ASN:O	2.21	0.41
1:C:3:LEU:HB2	1:C:128:VAL:HG11	2.04	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	319/330 (97%)	313 (98%)	6 (2%)	0	100	100
1	B	312/330 (94%)	305 (98%)	7 (2%)	0	100	100
1	C	300/330 (91%)	293 (98%)	7 (2%)	0	100	100
1	D	306/330 (93%)	302 (99%)	4 (1%)	0	100	100
All	All	1237/1320 (94%)	1213 (98%)	24 (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	272/277 (98%)	268 (98%)	4 (2%)	76	89
1	B	267/277 (96%)	259 (97%)	8 (3%)	53	70
1	C	260/277 (94%)	252 (97%)	8 (3%)	52	68
1	D	264/277 (95%)	257 (97%)	7 (3%)	57	74
All	All	1063/1108 (96%)	1036 (98%)	27 (2%)	60	77

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	LEU
1	A	18	THR
1	A	175	ASN
1	A	183	LYS
1	B	5	LEU
1	B	124	LEU
1	B	137	LEU
1	B	175	ASN
1	B	226	GLN
1	B	243	THR
1	B	269	ASN
1	B	272	ASP
1	C	18[A]	THR
1	C	18[B]	THR
1	C	22	LEU
1	C	24	ASP
1	C	129	GLU
1	C	175	ASN
1	C	183	LYS
1	C	187	SER
1	D	5	LEU
1	D	125	LEU
1	D	175	ASN
1	D	210	GLN
1	D	226	GLN
1	D	243	THR

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Mol	Chain	Res	Type
1	D	252	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

There are no ligands in this entry.

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	318/330 (96%)	0.10	7 (2%) 59 69	36, 42, 49, 53	0
1	B	312/330 (94%)	0.29	24 (7%) 13 20	35, 43, 51, 55	0
1	C	303/330 (91%)	0.41	19 (6%) 19 28	36, 43, 49, 55	0
1	D	310/330 (93%)	0.05	9 (2%) 49 59	35, 42, 49, 54	0
All	All	1243/1320 (94%)	0.21	59 (4%) 30 40	35, 42, 50, 55	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	250	VAL	6.4
1	A	250	VAL	6.2
1	B	247	LEU	5.4
1	B	290	THR	5.0
1	B	251	GLY	5.0
1	D	249	PRO	4.9
1	C	243	THR	4.9
1	C	292	TYR	4.6
1	C	250	VAL	4.4
1	D	288	ALA	4.3
1	C	126	GLU	3.9
1	B	245	SER	3.9
1	A	290	THR	3.8
1	A	288	ALA	3.7
1	D	194	ASN	3.6
1	D	251	GLY	3.6
1	B	306	GLU	3.6
1	C	154	GLN	3.4
1	B	250	VAL	3.4
1	C	267	HIS	3.3
1	B	288	ALA	3.2

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Mol	Chain	Res	Type	RSRZ
1	C	244	ILE	3.2
1	C	115	ALA	3.1
1	A	226	GLN	3.0
1	B	310	VAL	3.0
1	B	30	GLU	2.9
1	C	177	TYR	2.8
1	A	289	GLN	2.7
1	B	300	ASP	2.6
1	C	296	ASN	2.6
1	B	203	SER	2.6
1	C	0	HIS	2.6
1	D	292	TYR	2.5
1	C	75	ASP	2.5
1	B	239	VAL	2.4
1	B	223	LEU	2.4
1	B	200	SER	2.4
1	C	84	LYS	2.3
1	B	271	HIS	2.3
1	B	305	ILE	2.3
1	A	308	LEU	2.3
1	C	30	GLU	2.3
1	B	186	ILE	2.3
1	B	283	LEU	2.3
1	B	249	PRO	2.2
1	B	284	ASN	2.2
1	B	234	HIS	2.1
1	A	249	PRO	2.1
1	B	308	LEU	2.1
1	C	189	LEU	2.1
1	C	301	LEU	2.1
1	C	234	HIS	2.1
1	C	167	ALA	2.1
1	D	-5	VAL	2.0
1	B	241	ILE	2.0
1	B	269	ASN	2.0
1	C	249	PRO	2.0
1	D	75[A]	ASP	2.0
1	D	226	GLN	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.



### 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

### 6.4 Ligands ⓘ

There are no ligands in this entry.

### 6.5 Other polymers ⓘ

There are no such residues in this entry.