



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:03 PM GMT

PDB ID : 4URQ  
Title : Crystal Structure of GGDEF domain (I site mutant) from T.maritima  
Authors : Deepthi, A.; Liew, C.W.; Liang, Z.X.; Swamianthan, K.; Lescar, J.  
Deposited on : 2014-07-01  
Resolution : 2.50 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](mailto: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.

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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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
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) : trunk26865

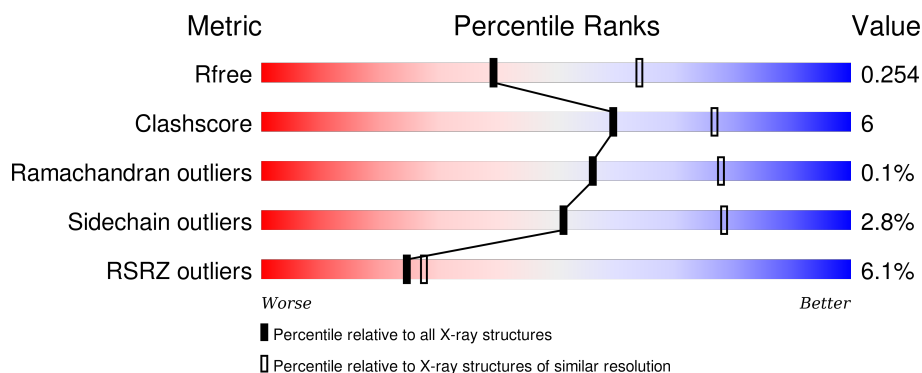
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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))
$R_{free}$	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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. 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	U	167	<div> <div>7%</div> <div> <div></div> <div>77%</div> <div>14%</div> <div>•</div> <div>8%</div> </div> </div>
1	V	167	<div> <div>5%</div> <div> <div></div> <div>75%</div> <div>16%</div> <div>•</div> <div>8%</div> </div> </div>
1	W	167	<div> <div>6%</div> <div> <div></div> <div>76%</div> <div>16%</div> <div></div> <div>8%</div> </div> </div>
1	X	167	<div> <div>7%</div> <div> <div></div> <div>76%</div> <div>16%</div> <div>•</div> <div>8%</div> </div> </div>
1	Y	167	<div> <div>5%</div> <div> <div></div> <div>81%</div> <div>10%</div> <div>•</div> <div>8%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	Z	167	<div><div></div><div>4%</div><div>77%</div><div>14%</div><div>8%</div></div>

## 2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 7658 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 DIGUANYLATE CYCLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	U	154	Total	C	N	O	S	0	0	0
			1258	809	216	229	4			
1	V	153	Total	C	N	O	S	0	0	0
			1250	803	215	228	4			
1	W	154	Total	C	N	O	S	0	1	0
			1257	810	213	229	5			
1	X	154	Total	C	N	O	S	0	0	0
			1258	809	216	229	4			
1	Y	154	Total	C	N	O	S	0	0	0
			1258	809	216	229	4			
1	Z	153	Total	C	N	O	S	0	0	0
			1250	803	215	228	4			

There are 60 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
U	90	MET	-	EXPRESSION TAG	UNP Q9X2A8
U	158	ALA	ARG	ENGINEERED MUTATION	UNP Q9X2A8
U	249	LEU	-	EXPRESSION TAG	UNP Q9X2A8
U	250	GLU	-	EXPRESSION TAG	UNP Q9X2A8
U	251	HIS	-	EXPRESSION TAG	UNP Q9X2A8
U	252	HIS	-	EXPRESSION TAG	UNP Q9X2A8
U	253	HIS	-	EXPRESSION TAG	UNP Q9X2A8
U	254	HIS	-	EXPRESSION TAG	UNP Q9X2A8
U	255	HIS	-	EXPRESSION TAG	UNP Q9X2A8
U	256	HIS	-	EXPRESSION TAG	UNP Q9X2A8
V	90	MET	-	EXPRESSION TAG	UNP Q9X2A8
V	158	ALA	ARG	ENGINEERED MUTATION	UNP Q9X2A8
V	249	LEU	-	EXPRESSION TAG	UNP Q9X2A8
V	250	GLU	-	EXPRESSION TAG	UNP Q9X2A8
V	251	HIS	-	EXPRESSION TAG	UNP Q9X2A8
V	252	HIS	-	EXPRESSION TAG	UNP Q9X2A8
V	253	HIS	-	EXPRESSION TAG	UNP Q9X2A8

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Chain	Residue	Modelled	Actual	Comment	Reference
V	254	HIS	-	EXPRESSION TAG	UNP Q9X2A8
V	255	HIS	-	EXPRESSION TAG	UNP Q9X2A8
V	256	HIS	-	EXPRESSION TAG	UNP Q9X2A8
W	90	MET	-	EXPRESSION TAG	UNP Q9X2A8
W	158	ALA	ARG	ENGINEERED MUTATION	UNP Q9X2A8
W	249	LEU	-	EXPRESSION TAG	UNP Q9X2A8
W	250	GLU	-	EXPRESSION TAG	UNP Q9X2A8
W	251	HIS	-	EXPRESSION TAG	UNP Q9X2A8
W	252	HIS	-	EXPRESSION TAG	UNP Q9X2A8
W	253	HIS	-	EXPRESSION TAG	UNP Q9X2A8
W	254	HIS	-	EXPRESSION TAG	UNP Q9X2A8
W	255	HIS	-	EXPRESSION TAG	UNP Q9X2A8
W	256	HIS	-	EXPRESSION TAG	UNP Q9X2A8
X	90	MET	-	EXPRESSION TAG	UNP Q9X2A8
X	158	ALA	ARG	ENGINEERED MUTATION	UNP Q9X2A8
X	249	LEU	-	EXPRESSION TAG	UNP Q9X2A8
X	250	GLU	-	EXPRESSION TAG	UNP Q9X2A8
X	251	HIS	-	EXPRESSION TAG	UNP Q9X2A8
X	252	HIS	-	EXPRESSION TAG	UNP Q9X2A8
X	253	HIS	-	EXPRESSION TAG	UNP Q9X2A8
X	254	HIS	-	EXPRESSION TAG	UNP Q9X2A8
X	255	HIS	-	EXPRESSION TAG	UNP Q9X2A8
X	256	HIS	-	EXPRESSION TAG	UNP Q9X2A8
Y	90	MET	-	EXPRESSION TAG	UNP Q9X2A8
Y	158	ALA	ARG	ENGINEERED MUTATION	UNP Q9X2A8
Y	249	LEU	-	EXPRESSION TAG	UNP Q9X2A8
Y	250	GLU	-	EXPRESSION TAG	UNP Q9X2A8
Y	251	HIS	-	EXPRESSION TAG	UNP Q9X2A8
Y	252	HIS	-	EXPRESSION TAG	UNP Q9X2A8
Y	253	HIS	-	EXPRESSION TAG	UNP Q9X2A8
Y	254	HIS	-	EXPRESSION TAG	UNP Q9X2A8
Y	255	HIS	-	EXPRESSION TAG	UNP Q9X2A8
Y	256	HIS	-	EXPRESSION TAG	UNP Q9X2A8
Z	90	MET	-	EXPRESSION TAG	UNP Q9X2A8
Z	158	ALA	ARG	ENGINEERED MUTATION	UNP Q9X2A8
Z	249	LEU	-	EXPRESSION TAG	UNP Q9X2A8
Z	250	GLU	-	EXPRESSION TAG	UNP Q9X2A8
Z	251	HIS	-	EXPRESSION TAG	UNP Q9X2A8
Z	252	HIS	-	EXPRESSION TAG	UNP Q9X2A8
Z	253	HIS	-	EXPRESSION TAG	UNP Q9X2A8
Z	254	HIS	-	EXPRESSION TAG	UNP Q9X2A8
Z	255	HIS	-	EXPRESSION TAG	UNP Q9X2A8

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Chain	Residue	Modelled	Actual	Comment	Reference
Z	256	HIS	-	EXPRESSION TAG	UNP Q9X2A8

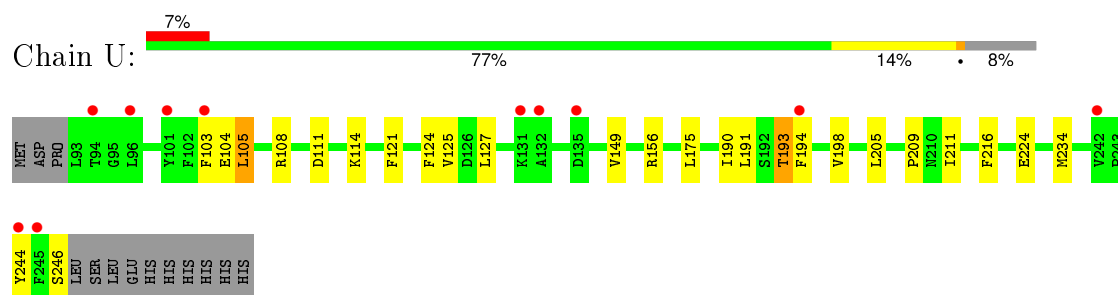
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	U	22	Total O 22 22	0	0
2	V	26	Total O 26 26	0	0
2	W	23	Total O 23 23	0	0
2	X	19	Total O 19 19	0	0
2	Y	18	Total O 18 18	0	0
2	Z	19	Total O 19 19	0	0

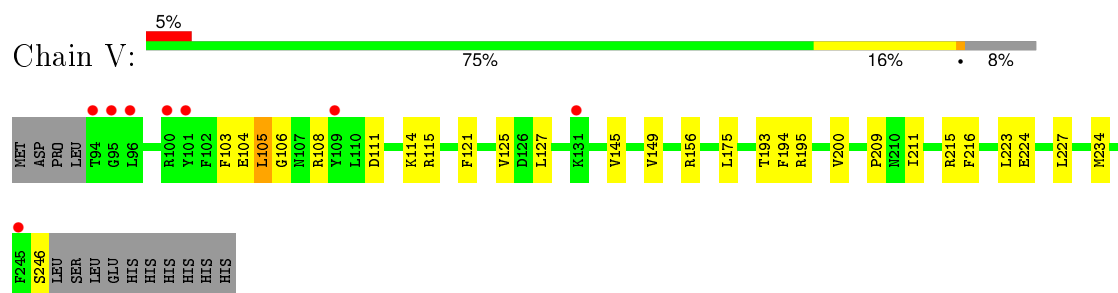
### 3 Residue-property plots [i](#)

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.

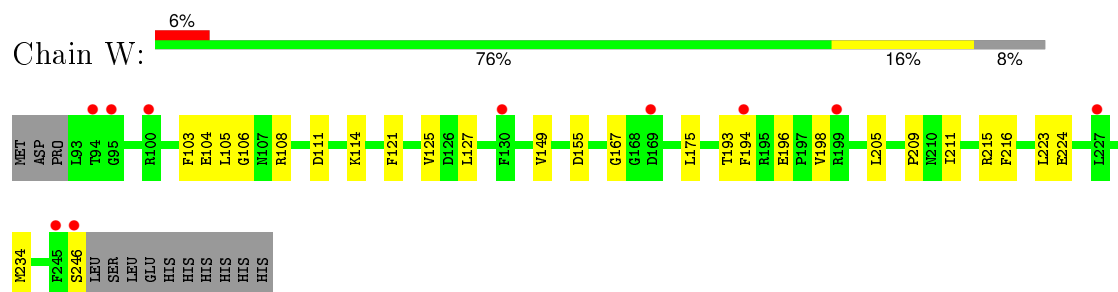
#### • Molecule 1: DIGUANYLATE CYCLASE



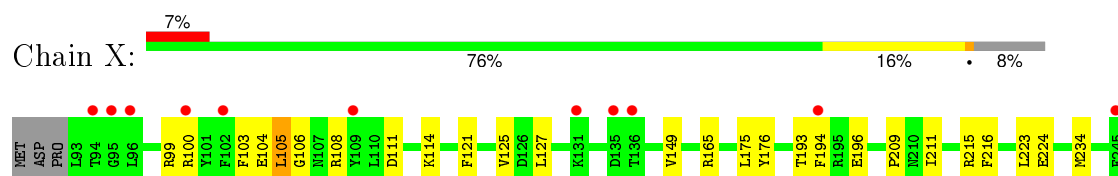
#### • Molecule 1: DIGUANYLATE CYCLASE



#### • Molecule 1: DIGUANYLATE CYCLASE

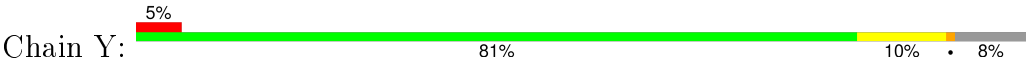


#### • Molecule 1: DIGUANYLATE CYCLASE



S246
LEU
SER
LEU
GLU
HIS
HIS
HIS
HIS
HIS

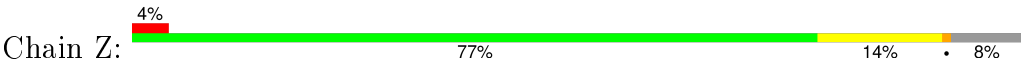
● Molecule 1: DIGUANYLATE CYCLASE



MET	ASP	PRO	L83	T94	G95	Y101	E104	L105	R108	Y109	L110	D111	K114	F121	V125	D135	L140	L146	L175	I190	T193	F194	R199	P209	N210	I211	R215	F216	M234	E239	F245	S246	LEU	SER	LEU	GLU	HIS	HIS	HIS	HIS	HIS
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HIS
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● Molecule 1: DIGUANYLATE CYCLASE



MET	ASP	PRO	T94	R100	Y101	F102	F103	E104	L105	G106	R107	R108	D111	K114	K119	V125	D126	L127	D135	V149	R165	G168	D169	T193	F194	R195	E196	P209	H210	I211	R215	F216	L223	E224	M234	Y244	F245	S246	LEU	SER	LEU	GLU	HIS
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## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	159.41Å 91.98Å 87.59Å 90.00° 90.04° 90.00°	Depositor
Resolution (Å)	38.37 – 2.50 36.28 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.7 (38.37-2.50) 99.4 (36.28-2.50)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.67 (at 2.51Å)	Xtriage
Refinement program	BUSTER 2.10.0	Depositor
R, $R_{free}$	0.222 , 0.247 0.228 , 0.254	Depositor DCC
$R_{free}$ test set	2196 reflections (5.29%)	DCC
Wilson B-factor (Å <sup>2</sup> )	45.5	Xtriage
Anisotropy	0.048	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 35.9	EDS
Estimated twinning fraction	0.479 for -1/2*h-3/2*k,-1/2*h+1/2*k,-l 0.478 for -1/2*h+3/2*k,1/2*h+1/2*k,-l 0.478 for 1/2*h-3/2*k,-1/2*h-1/2*k,-l 0.479 for 1/2*h+3/2*k,1/2*h-1/2*k,-l 0.487 for -h,-k,l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 43743 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	7658	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.92% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<sup>2</sup> Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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	U	0.51	0/1281	0.72	0/1721
1	V	0.52	0/1273	0.70	0/1710
1	W	0.52	0/1283	0.71	0/1724
1	X	0.51	0/1281	0.71	0/1721
1	Y	0.51	0/1281	0.71	0/1721
1	Z	0.51	0/1273	0.70	0/1710
All	All	0.51	0/7672	0.71	0/10307

There are no bond length outliers.

There are no bond angle outliers.

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	U	1258	0	1283	17	0
1	V	1250	0	1272	20	0
1	W	1257	0	1281	17	0
1	X	1258	0	1283	17	0
1	Y	1258	0	1283	13	0
1	Z	1250	0	1272	14	0
2	U	22	0	0	0	0
2	V	26	0	0	0	0
2	W	23	0	0	0	0
2	X	19	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	Y	18	0	0	0	0
2	Z	19	0	0	0	0
All	All	7658	0	7674	87	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 6.

All (87) 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:X:215:ARG:HH22	1:X:246:SER:HB2	1.27	1.00
1:Y:215:ARG:HH22	1:Y:246:SER:HB2	1.27	1.00
1:V:215:ARG:HH22	1:V:246:SER:HB2	1.28	0.96
1:Z:215:ARG:HH22	1:Z:246:SER:HB2	1.28	0.96
1:W:215:ARG:HH22	1:W:246:SER:HB2	1.27	0.96
1:V:127:LEU:HD11	1:V:149:VAL:HG11	1.74	0.69
1:V:194:PHE:HB2	1:V:209:PRO:HG3	1.76	0.68
1:U:194:PHE:HB2	1:U:209:PRO:HG3	1.76	0.67
1:Z:194:PHE:HB2	1:Z:209:PRO:HG3	1.75	0.67
1:Y:194:PHE:HB2	1:Y:209:PRO:HG3	1.76	0.67
1:W:194:PHE:HB2	1:W:209:PRO:HG3	1.76	0.66
1:V:108:ARG:HG3	1:X:105:LEU:HD21	1.79	0.65
1:V:103:PHE:HE2	1:V:224:GLU:HG3	1.61	0.65
1:W:103:PHE:HE2	1:W:224:GLU:HG3	1.61	0.64
1:X:215:ARG:HH22	1:X:246:SER:CB	2.09	0.64
1:U:105:LEU:HD21	1:Y:108:ARG:HG3	1.80	0.64
1:Z:103:PHE:HE2	1:Z:224:GLU:HG3	1.61	0.64
1:X:103:PHE:HE2	1:X:224:GLU:HG3	1.61	0.64
1:U:103:PHE:HE2	1:U:224:GLU:HG3	1.61	0.64
1:U:108:ARG:HG3	1:Y:105:LEU:HD21	1.80	0.64
1:V:105:LEU:HD21	1:X:108:ARG:HG3	1.82	0.62
1:Y:125:VAL:HG22	1:Y:211:ILE:HG12	1.79	0.62
1:V:114:LYS:HG3	1:V:216:PHE:HZ	1.68	0.59
1:U:114:LYS:HG3	1:U:216:PHE:HZ	1.68	0.58
1:W:114:LYS:HG3	1:W:216:PHE:HZ	1.68	0.58
1:W:125:VAL:HG22	1:W:211:ILE:HG12	1.85	0.58
1:V:125:VAL:HG22	1:V:211:ILE:HG12	1.85	0.57
1:V:215:ARG:HH22	1:V:246:SER:CB	2.09	0.57
1:X:194:PHE:HB2	1:X:209:PRO:HG3	1.87	0.57
1:X:125:VAL:HG22	1:X:211:ILE:HG12	1.85	0.57
1:U:125:VAL:HG22	1:U:211:ILE:HG12	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Z:114:LYS:HG3	1:Z:216:PHE:HZ	1.69	0.57
1:X:127:LEU:HD11	1:X:149:VAL:HG11	1.88	0.56
1:Y:114:LYS:HG3	1:Y:216:PHE:HZ	1.69	0.56
1:X:114:LYS:HG3	1:X:216:PHE:HZ	1.69	0.56
1:Z:215:ARG:HH22	1:Z:246:SER:CB	2.09	0.56
1:W:215:ARG:HH22	1:W:246:SER:CB	2.09	0.56
1:Z:125:VAL:HG22	1:Z:211:ILE:HG12	1.87	0.56
1:Y:215:ARG:HH22	1:Y:246:SER:CB	2.09	0.54
1:U:156:ARG:HH22	1:Z:196:GLU:HG2	1.74	0.53
1:V:156:ARG:HH22	1:W:196:GLU:HG2	1.75	0.51
1:W:114:LYS:HG3	1:W:216:PHE:CZ	2.47	0.48
1:Y:114:LYS:HG3	1:Y:216:PHE:CZ	2.48	0.48
1:Y:190:ILE:O	1:Y:193:THR:HG22	2.14	0.48
1:X:114:LYS:HG3	1:X:216:PHE:CZ	2.49	0.48
1:U:191:LEU:HD11	1:U:244:TYR:HD2	1.79	0.48
1:Z:114:LYS:HG3	1:Z:216:PHE:CZ	2.49	0.47
1:Y:111:ASP:HA	1:Y:114:LYS:HD2	1.96	0.47
1:Z:127:LEU:CD1	1:Z:149:VAL:HG11	2.43	0.47
1:V:145:VAL:HG23	1:V:200:VAL:HG21	1.97	0.47
1:V:114:LYS:HG3	1:V:216:PHE:CZ	2.48	0.47
1:Z:111:ASP:HA	1:Z:114:LYS:HD2	1.96	0.47
1:V:156:ARG:NH2	1:W:196:GLU:HG2	2.29	0.47
1:W:111:ASP:HA	1:W:114:LYS:HD2	1.96	0.47
1:U:114:LYS:HG3	1:U:216:PHE:CZ	2.48	0.47
1:X:111:ASP:HA	1:X:114:LYS:HD2	1.96	0.47
1:W:127:LEU:HD11	1:W:149:VAL:HG11	1.97	0.47
1:U:111:ASP:HA	1:U:114:LYS:HD2	1.96	0.46
1:V:111:ASP:HA	1:V:114:LYS:HD2	1.96	0.46
1:U:127:LEU:CD1	1:U:149:VAL:HG11	2.46	0.45
1:X:99:ARG:HE	1:X:100:ARG:NH2	2.15	0.44
1:U:190:ILE:O	1:U:193:THR:HG22	2.18	0.44
1:U:104:GLU:O	1:U:108:ARG:HG2	2.18	0.44
1:V:104:GLU:O	1:V:108:ARG:HG2	2.17	0.43
1:X:104:GLU:O	1:X:108:ARG:HG2	2.18	0.43
1:V:115:ARG:NH2	1:X:176:TYR:O	2.43	0.43
1:Z:104:GLU:O	1:Z:108:ARG:HG2	2.18	0.43
1:W:104:GLU:O	1:W:108:ARG:HG2	2.19	0.43
1:Y:104:GLU:O	1:Y:108:ARG:HG2	2.19	0.42
1:W:127:LEU:CD1	1:W:149:VAL:HG11	2.50	0.42
1:X:100:ARG:HA	1:X:100:ARG:HD3	1.92	0.42
1:Z:165:ARG:NH1	1:Z:168:GLY:O	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:106:GLY:HA3	1:X:223:LEU:HD22	2.02	0.42
1:W:106:GLY:HA3	1:W:223:LEU:HD22	2.02	0.41
1:U:198:VAL:N	1:U:205:LEU:O	2.52	0.41
1:Z:106:GLY:HA3	1:Z:223:LEU:HD22	2.03	0.41
1:W:198:VAL:N	1:W:205:LEU:O	2.54	0.41
1:U:156:ARG:NH2	1:Z:196:GLU:HG2	2.34	0.41
1:V:106:GLY:HA3	1:V:223:LEU:HD22	2.02	0.41
1:V:227:LEU:HB3	1:Y:140:LEU:HD11	2.02	0.40
1:U:124:PHE:O	1:U:211:ILE:HA	2.21	0.40
1:W:121:PHE:HB2	1:W:175:LEU:HB2	2.03	0.40
1:Y:121:PHE:HB2	1:Y:175:LEU:HB2	2.03	0.40
1:U:121:PHE:HB2	1:U:175:LEU:HB2	2.03	0.40
1:V:121:PHE:HB2	1:V:175:LEU:HB2	2.03	0.40
1:X:121:PHE:HB2	1:X:175:LEU:HB2	2.03	0.40
1:V:195:ARG:NH1	1:W:155:ASP:O	2.54	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	U	152/167 (91%)	147 (97%)	5 (3%)	0	100	100
1	V	151/167 (90%)	146 (97%)	5 (3%)	0	100	100
1	W	153/167 (92%)	145 (95%)	7 (5%)	1 (1%)	26	46
1	X	152/167 (91%)	145 (95%)	7 (5%)	0	100	100
1	Y	152/167 (91%)	145 (95%)	7 (5%)	0	100	100
1	Z	151/167 (90%)	144 (95%)	7 (5%)	0	100	100
All	All	911/1002 (91%)	872 (96%)	38 (4%)	1 (0%)	56	78

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	W	167	GLY

### 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	U	136/149 (91%)	132 (97%)	4 (3%)	50	77
1	V	135/149 (91%)	132 (98%)	3 (2%)	60	84
1	W	136/149 (91%)	133 (98%)	3 (2%)	60	84
1	X	136/149 (91%)	131 (96%)	5 (4%)	41	68
1	Y	136/149 (91%)	133 (98%)	3 (2%)	60	84
1	Z	135/149 (91%)	130 (96%)	5 (4%)	41	68
All	All	814/894 (91%)	791 (97%)	23 (3%)	51	78

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	U	105	LEU
1	U	193	THR
1	U	234	MET
1	U	246	SER
1	V	105	LEU
1	V	193	THR
1	V	234	MET
1	W	105	LEU
1	W	193	THR
1	W	234	MET
1	X	105	LEU
1	X	165	ARG
1	X	193	THR
1	X	196	GLU
1	X	234	MET
1	Y	105	LEU
1	Y	193	THR
1	Y	234	MET

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Mol	Chain	Res	Type
1	Z	105	LEU
1	Z	119	LYS
1	Z	165	ARG
1	Z	193	THR
1	Z	234	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	U	222	ASN
1	V	222	ASN
1	W	222	ASN
1	X	134	ASN
1	X	222	ASN
1	Z	222	ASN

### 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](#)

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 ⓘ

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	U	154/167 (92%)	0.63	11 (7%)	19 21	30, 51, 83, 98	0
1	V	153/167 (91%)	0.60	8 (5%)	31 35	30, 51, 82, 96	0
1	W	154/167 (92%)	0.64	10 (6%)	22 25	31, 50, 84, 101	0
1	X	154/167 (92%)	0.72	11 (7%)	19 21	30, 51, 83, 105	0
1	Y	154/167 (92%)	0.62	9 (5%)	26 30	30, 50, 85, 104	0
1	Z	153/167 (91%)	0.51	7 (4%)	36 41	31, 50, 83, 96	0
All	All	922/1002 (92%)	0.62	56 (6%)	25 27	30, 51, 84, 105	0

All (56) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	V	101	TYR	4.5
1	V	94	THR	4.4
1	X	131	LYS	4.3
1	W	94	THR	3.8
1	U	96	LEU	3.8
1	Y	199	ARG	3.7
1	X	96	LEU	3.7
1	Y	95	GLY	3.6
1	X	245	PHE	3.6
1	W	169	ASP	3.6
1	Y	101	TYR	3.5
1	U	131	LYS	3.4
1	Z	135	ASP	3.1
1	W	245	PHE	3.1
1	Y	245	PHE	3.1
1	Z	244	TYR	3.0
1	V	245	PHE	3.0
1	U	245	PHE	2.9
1	V	109	TYR	2.9

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Mol	Chain	Res	Type	RSRZ
1	W	100	ARG	2.8
1	V	96	LEU	2.8
1	Z	245	PHE	2.7
1	V	100	ARG	2.7
1	V	131	LYS	2.7
1	X	135	ASP	2.6
1	X	95	GLY	2.6
1	Z	101	TYR	2.6
1	U	101	TYR	2.6
1	V	95	GLY	2.5
1	X	100	ARG	2.5
1	W	130	PHE	2.5
1	Y	135	ASP	2.4
1	U	194	PHE	2.4
1	Z	100	ARG	2.3
1	Y	121	PHE	2.3
1	X	94	THR	2.3
1	U	135	ASP	2.3
1	X	102	PHE	2.3
1	Y	109	TYR	2.3
1	U	132	ALA	2.2
1	Z	196	GLU	2.2
1	W	199	ARG	2.2
1	Z	169	ASP	2.2
1	W	95	GLY	2.2
1	U	242	VAL	2.2
1	Y	239	GLU	2.2
1	X	109	TYR	2.2
1	U	244	TYR	2.1
1	X	194	PHE	2.1
1	Y	146	LEU	2.1
1	U	94	THR	2.1
1	W	194	PHE	2.1
1	X	136	THR	2.1
1	U	103	PHE	2.0
1	W	227	LEU	2.0
1	W	246	SER	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 [i](#)

There are no carbohydrates 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.