



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 12:09 AM UTC

PDB ID : 5DEC / pdb\_00005dec  
Title : Crystal structure of the small alarmone synthetase 1 from *Bacillus subtilis*  
Authors : Steinchen, W.; Altegoer, A.; Schuhmacher, J.S.; Bange, G.  
Deposited on : 2015-08-25  
Resolution : 2.00 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

<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  
Xtrriage (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

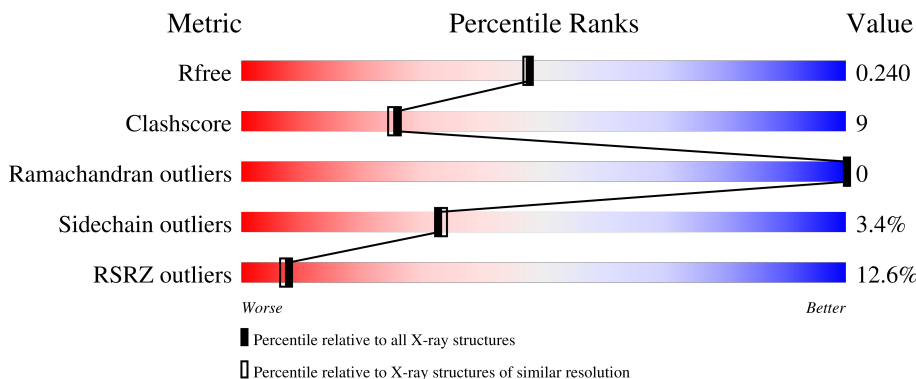
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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  $\geq 3$ , 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	218	 9% 72% 14% 14%
1	B	218	 12% 71% 17% 12%
1	C	218	 10% 65% 14% 5% 16%
1	D	218	 13% 67% 18% 13%

## 2 Entry composition i

There are 2 unique types of molecules in this entry. The entry contains 6542 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 GTP pyrophosphokinase YjbM.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	188	Total 1546	C 985	N 270	O 285	S 6	0	0	0
1	B	191	Total 1572	C 1001	N 276	O 289	S 6	0	0	0
1	C	183	Total 1507	C 960	N 262	O 279	S 6	0	0	0
1	D	189	Total 1552	C 988	N 271	O 287	S 6	0	0	0

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	MET	-	expression tag	UNP O31611
A	-5	GLY	-	expression tag	UNP O31611
A	-4	HIS	-	expression tag	UNP O31611
A	-3	HIS	-	expression tag	UNP O31611
A	-2	HIS	-	expression tag	UNP O31611
A	-1	HIS	-	expression tag	UNP O31611
A	0	HIS	-	expression tag	UNP O31611
A	1	HIS	-	expression tag	UNP O31611
A	4	ASP	LYS	conflict	UNP O31611
A	5	LYS	GLN	conflict	UNP O31611
B	-6	MET	-	expression tag	UNP O31611
B	-5	GLY	-	expression tag	UNP O31611
B	-4	HIS	-	expression tag	UNP O31611
B	-3	HIS	-	expression tag	UNP O31611
B	-2	HIS	-	expression tag	UNP O31611
B	-1	HIS	-	expression tag	UNP O31611
B	0	HIS	-	expression tag	UNP O31611
B	1	HIS	-	expression tag	UNP O31611
B	4	ASP	LYS	conflict	UNP O31611
B	5	LYS	GLN	conflict	UNP O31611
C	-6	MET	-	expression tag	UNP O31611

*Continued on next page...*

*Continued from previous page...*

Chain	Residue	Modelled	Actual	Comment	Reference
C	-5	GLY	-	expression tag	UNP O31611
C	-4	HIS	-	expression tag	UNP O31611
C	-3	HIS	-	expression tag	UNP O31611
C	-2	HIS	-	expression tag	UNP O31611
C	-1	HIS	-	expression tag	UNP O31611
C	0	HIS	-	expression tag	UNP O31611
C	1	HIS	-	expression tag	UNP O31611
C	4	ASP	LYS	conflict	UNP O31611
C	5	LYS	GLN	conflict	UNP O31611
D	-6	MET	-	expression tag	UNP O31611
D	-5	GLY	-	expression tag	UNP O31611
D	-4	HIS	-	expression tag	UNP O31611
D	-3	HIS	-	expression tag	UNP O31611
D	-2	HIS	-	expression tag	UNP O31611
D	-1	HIS	-	expression tag	UNP O31611
D	0	HIS	-	expression tag	UNP O31611
D	1	HIS	-	expression tag	UNP O31611
D	4	ASP	LYS	conflict	UNP O31611
D	5	LYS	GLN	conflict	UNP O31611

- Molecule 2 is water.

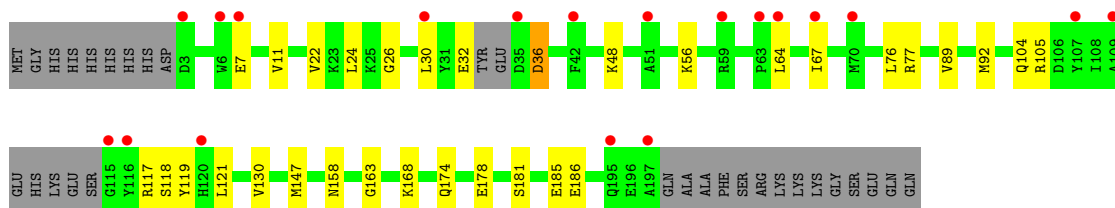
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	90	Total O 90 90	0	0
2	B	96	Total O 96 96	0	0
2	C	92	Total O 92 92	0	0
2	D	87	Total O 87 87	0	0

### 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.

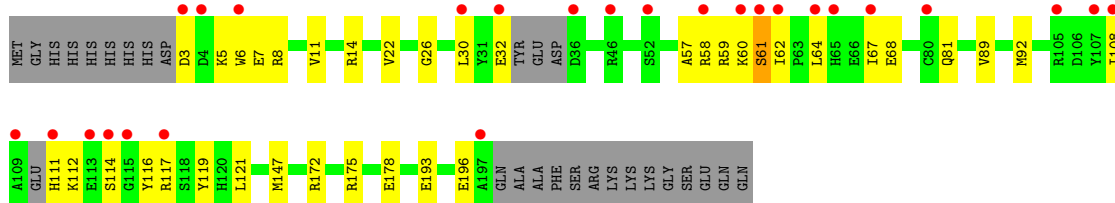
- Molecule 1: GTP pyrophosphokinase YjbM

Chain A: 



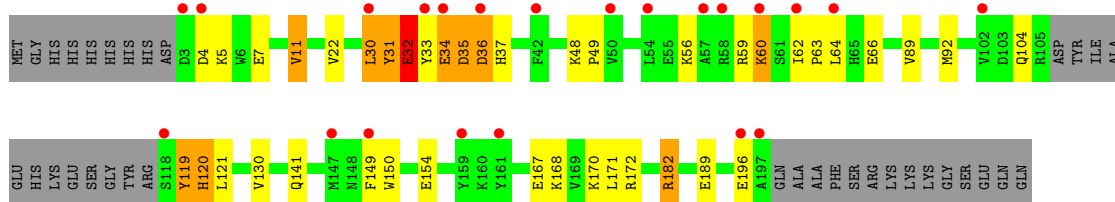
- Molecule 1: GTP pyrophosphokinase YjbM

Chain B: 



- Molecule 1: GTP pyrophosphokinase YjbM

Chain C: 



- Molecule 1: GTP pyrophosphokinase YjbM

Chain D: 



## 4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	78.13Å 77.69Å 81.61Å 90.00° 90.40° 90.00°	Depositor
Resolution (Å)	56.63 – 2.00 56.63 – 2.00	Depositor EDS
% Data completeness (in resolution range)	97.4 (56.63-2.00) 96.3 (56.63-2.00)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.79 (at 2.00Å)	Xtriage
Refinement program	PHENIX 1.9_1685	Depositor
R, $R_{free}$	0.194 , 0.238 0.199 , 0.240	Depositor DCC
$R_{free}$ test set	2025 reflections (3.06%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	32.0	Xtriage
Anisotropy	0.248	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 50.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.000 for l,k,-h 0.017 for -h,-l,-k 0.001 for -h,l,k 0.023 for k,h,-l 0.034 for -k,-h,-l 0.002 for l,h,k 0.002 for k,l,h 0.001 for -l,-h,k 0.001 for -k,-l,h 0.026 for h,-k,-l 0.016 for l,-k,h	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	6542	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.79% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality i

### 5.1 Standard geometry i

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	0.58	0/1572	0.84	1/2116 (0.0%)
1	B	0.54	0/1599	0.82	1/2151 (0.0%)
1	C	0.66	0/1533	0.97	8/2065 (0.4%)
1	D	0.61	0/1578	0.97	4/2124 (0.2%)
All	All	0.60	0/6282	0.90	14/8456 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1
1	D	0	1
All	All	0	2

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	59	ARG	N-CA-C	-12.52	98.01	113.28
1	C	34	GLU	N-CA-C	10.86	123.11	111.28
1	C	31	TYR	N-CA-C	8.47	120.59	111.36
1	C	4	ASP	N-CA-C	-7.46	103.64	112.89
1	D	58	ARG	CA-CB-CG	6.16	126.42	114.10
1	C	62	ILE	CA-C-N	6.11	126.05	119.76
1	C	62	ILE	C-N-CA	6.11	126.05	119.76
1	D	108	ILE	N-CA-C	-5.88	104.59	111.00
1	D	61	SER	N-CA-C	5.74	119.32	111.39
1	A	36	ASP	N-CA-C	5.48	117.84	109.95
1	C	11	VAL	CB-CA-C	-5.22	108.74	113.70
1	C	119	TYR	CA-C-N	-5.07	114.85	122.81

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	119	TYR	C-N-CA	-5.07	114.85	122.81
1	B	5	LYS	N-CA-C	-5.04	105.91	111.71

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	32	GLU	Peptide
1	D	58	ARG	Peptide

## 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1546	0	1567	26	0
1	B	1572	0	1594	31	1
1	C	1507	0	1529	39	0
1	D	1552	0	1572	33	1
2	A	90	0	0	2	0
2	B	96	0	0	4	0
2	C	92	0	0	5	0
2	D	87	0	0	9	0
All	All	6542	0	6262	118	1

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 9.

All (118) 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:C:33:TYR:C	1:C:35:ASP:OD1	1.79	1.25
1:D:32:GLU:C	1:D:35:ASP:OD2	1.80	1.24
1:D:32:GLU:O	1:D:35:ASP:CG	1.84	1.18
1:C:32:GLU:C	1:C:35:ASP:OD1	1.93	1.10
1:C:32:GLU:O	1:C:35:ASP:OD2	1.85	0.94

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:22:VAL:HG21	1:C:22:VAL:HG21	1.48	0.94
1:A:158:ASN:ND2	1:A:163:GLY:HA2	1.82	0.93
1:C:36:ASP:OD1	1:C:37:HIS:N	2.03	0.91
1:D:147:MET:SD	2:D:352:HOH:O	2.30	0.89
1:C:167:GLU:O	2:C:301:HOH:O	1.90	0.88
1:B:22:VAL:HG21	1:D:22:VAL:HG21	1.55	0.88
1:C:32:GLU:O	1:C:35:ASP:CG	2.18	0.86
1:C:33:TYR:CA	1:C:35:ASP:OD1	2.24	0.84
1:D:192:GLY:O	2:D:301:HOH:O	1.94	0.84
1:C:34:GLU:N	1:C:35:ASP:OD1	2.10	0.82
1:B:147:MET:SD	2:B:383:HOH:O	2.38	0.81
1:C:5:LYS:NZ	2:C:302:HOH:O	2.14	0.81
1:A:158:ASN:HD21	1:A:163:GLY:HA2	1.46	0.80
1:D:196:GLU:N	2:D:301:HOH:O	2.14	0.79
1:D:168:LYS:O	2:D:302:HOH:O	2.02	0.77
1:C:32:GLU:C	1:C:35:ASP:CG	2.52	0.76
1:C:32:GLU:O	1:C:35:ASP:OD1	2.03	0.76
1:B:196:GLU:OE2	1:C:168:LYS:NZ	2.18	0.76
1:C:31:TYR:O	1:C:34:GLU:HB2	1.88	0.73
1:D:55:GLU:O	1:D:59:ARG:HG2	1.90	0.72
1:B:193:GLU:OE2	1:C:172:ARG:NH1	2.24	0.71
1:C:33:TYR:N	1:C:35:ASP:OD1	2.23	0.71
1:B:61:SER:O	1:B:61:SER:OG	2.05	0.70
1:B:32:GLU:O	2:B:301:HOH:O	2.10	0.70
1:A:130:VAL:HG23	1:C:30:LEU:HD23	1.72	0.70
1:A:119:TYR:HE2	1:A:121:LEU:HD23	1.58	0.69
1:B:3:ASP:N	2:B:303:HOH:O	2.27	0.68
1:B:178:GLU:OE1	2:B:302:HOH:O	2.11	0.67
1:B:7:GLU:OE2	1:B:14:ARG:NH2	2.28	0.66
1:A:186:GLU:HG3	2:D:303:HOH:O	1.96	0.65
1:B:119:TYR:HE2	1:B:121:LEU:HD23	1.62	0.64
1:B:193:GLU:O	1:B:196:GLU:HG2	1.98	0.63
1:C:56:LYS:NZ	1:C:60:LYS:HD2	2.14	0.61
1:B:112:LYS:HE2	1:B:114:SER:OG	2.00	0.61
1:D:58:ARG:HB2	1:D:62:ILE:H	1.65	0.60
1:D:182:ARG:NH1	2:D:306:HOH:O	2.25	0.59
1:C:120:HIS:ND1	1:C:141:GLN:HG2	2.18	0.59
1:D:119:TYR:HE2	1:D:121:LEU:HD23	1.66	0.58
1:D:13:TYR:OH	1:D:68:GLU:HG3	2.04	0.58
1:C:170:LYS:N	2:C:301:HOH:O	2.37	0.57
1:D:32:GLU:O	1:D:35:ASP:OD2	0.63	0.57

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:117:ARG:C	1:D:147:MET:HE3	2.30	0.57
1:A:105:ARG:HH11	1:A:105:ARG:HG3	1.70	0.56
1:C:171:LEU:N	2:C:301:HOH:O	1.95	0.56
1:D:58:ARG:HD2	1:D:61:SER:HA	1.88	0.55
1:D:7:GLU:O	1:D:11:VAL:HG23	2.07	0.54
1:D:175:ARG:O	2:D:303:HOH:O	2.18	0.53
1:B:81:GLN:HG3	1:C:149:PHE:HE1	1.74	0.53
1:A:117:ARG:C	1:A:147:MET:HE3	2.34	0.52
1:B:64:LEU:H	1:B:64:LEU:HD12	1.74	0.52
1:D:64:LEU:HA	1:D:67:ILE:HG23	1.91	0.52
1:A:48:LYS:HZ2	1:A:56:LYS:HD3	1.75	0.52
1:A:89:VAL:HA	1:A:92:MET:HE2	1.91	0.52
1:A:7:GLU:O	1:A:11:VAL:HG23	2.11	0.51
1:C:56:LYS:HZ3	1:C:60:LYS:HD2	1.74	0.51
1:D:55:GLU:HB3	1:D:59:ARG:HE	1.74	0.51
1:B:117:ARG:C	1:B:147:MET:HE3	2.37	0.50
1:C:7:GLU:O	1:C:11:VAL:HG23	2.12	0.50
1:D:165:ILE:HG12	2:D:325:HOH:O	2.11	0.50
1:D:157:LEU:HB3	1:D:165:ILE:HD12	1.94	0.49
1:C:63:PRO:HD2	1:C:66:GLU:OE1	2.11	0.49
1:B:26:GLY:O	1:B:30:LEU:HD13	2.13	0.49
1:A:174:GLN:NE2	1:A:178:GLU:OE2	2.32	0.49
1:B:111:HIS:HB2	1:B:116:TYR:O	2.14	0.48
1:C:168:LYS:C	2:C:301:HOH:O	2.55	0.48
1:C:150:TRP:CH2	1:C:154:GLU:OE1	2.66	0.48
1:C:104:GLN:OE1	1:C:119:TYR:OH	2.19	0.48
1:A:30:LEU:HD23	1:C:130:VAL:HG22	1.96	0.47
1:B:3:ASP:O	1:B:6:TRP:HB2	2.14	0.47
1:A:77:ARG:NH1	2:A:301:HOH:O	2.20	0.47
1:C:119:TYR:HE2	1:C:121:LEU:HD23	1.79	0.47
1:A:89:VAL:HA	1:A:92:MET:CE	2.45	0.47
1:C:36:ASP:CG	1:C:37:HIS:N	2.72	0.47
1:B:81:GLN:HG3	1:C:149:PHE:CE1	2.49	0.47
1:B:59:ARG:CZ	1:B:60:LYS:HE2	2.45	0.46
1:B:57:ALA:HB1	1:B:62:ILE:HB	1.96	0.46
1:B:7:GLU:O	1:B:11:VAL:HG23	2.15	0.46
1:D:58:ARG:HB2	1:D:62:ILE:N	2.30	0.46
1:D:58:ARG:HB2	1:D:61:SER:CA	2.46	0.45
1:A:26:GLY:O	1:A:30:LEU:HD13	2.17	0.45
1:C:120:HIS:ND1	1:C:141:GLN:CG	2.79	0.45
1:A:118:SER:HA	1:A:147:MET:CE	2.47	0.45

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:89:VAL:HA	1:B:92:MET:CE	2.47	0.45
1:B:59:ARG:C	1:B:60:LYS:HD3	2.42	0.45
1:C:89:VAL:HA	1:C:92:MET:CE	2.48	0.44
1:D:24:LEU:HD13	1:D:78:ILE:HD11	1.99	0.44
1:C:182:ARG:HA	1:C:182:ARG:HD3	1.80	0.44
1:D:58:ARG:HB2	1:D:61:SER:HA	1.99	0.44
1:A:168:LYS:HG2	1:D:193:GLU:CD	2.43	0.43
1:D:179:ALA:HB2	2:D:303:HOH:O	2.17	0.43
1:B:172:ARG:HE	1:C:189:GLU:HG2	1.82	0.43
1:A:105:ARG:HG3	1:A:105:ARG:NH1	2.32	0.43
1:A:104:GLN:HA	1:A:121:LEU:HA	2.00	0.43
1:A:181:SER:O	1:A:185:GLU:HG2	2.19	0.43
1:C:48:LYS:HA	1:C:49:PRO:HD3	1.91	0.42
1:B:196:GLU:HG3	1:B:196:GLU:O	2.19	0.42
1:D:57:ALA:O	1:D:62:ILE:HB	2.19	0.42
1:B:64:LEU:HA	1:B:67:ILE:HG23	2.02	0.42
1:C:104:GLN:HA	1:C:121:LEU:HA	2.01	0.42
1:D:53:ILE:O	1:D:57:ALA:N	2.52	0.42
1:D:178:GLU:HB3	1:D:182:ARG:NH2	2.34	0.42
1:A:24:LEU:HD12	1:A:76:LEU:HD22	2.01	0.42
1:D:26:GLY:O	1:D:30:LEU:HG	2.19	0.42
1:A:130:VAL:HG23	1:C:30:LEU:CD2	2.47	0.42
1:B:112:LYS:HE2	1:B:114:SER:HG	1.85	0.41
1:D:89:VAL:HA	1:D:92:MET:CE	2.49	0.41
1:B:108:ILE:HD11	1:B:147:MET:HE1	2.02	0.41
1:A:117:ARG:HD3	2:A:308:HOH:O	2.21	0.41
1:A:117:ARG:HB2	1:A:147:MET:HG2	2.03	0.41
1:B:64:LEU:HD12	1:B:64:LEU:N	2.36	0.41
1:D:117:ARG:O	1:D:147:MET:HE3	2.21	0.41
1:B:64:LEU:H	1:B:64:LEU:CD1	2.35	0.40
1:A:64:LEU:HA	1:A:67:ILE:HG23	2.04	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:175:ARG:NH2	1:D:186:GLU:OE1[1_556]	2.13	0.07

## 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	182/218 (84%)	180 (99%)	2 (1%)	0	100	100
1	B	185/218 (85%)	181 (98%)	4 (2%)	0	100	100
1	C	179/218 (82%)	178 (99%)	1 (1%)	0	100	100
1	D	183/218 (84%)	179 (98%)	4 (2%)	0	100	100
All	All	729/872 (84%)	718 (98%)	11 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	169/195 (87%)	167 (99%)	2 (1%)	63	70
1	B	172/195 (88%)	168 (98%)	4 (2%)	44	49
1	C	166/195 (85%)	156 (94%)	10 (6%)	17	14
1	D	170/195 (87%)	163 (96%)	7 (4%)	27	26
All	All	677/780 (87%)	654 (97%)	23 (3%)	32	33

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

Mol	Chain	Res	Type
1	A	32	GLU
1	A	36	ASP

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	B	8	ARG
1	B	58	ARG
1	B	61	SER
1	B	68	GLU
1	C	30	LEU
1	C	32	GLU
1	C	35	ASP
1	C	36	ASP
1	C	59	ARG
1	C	60	LYS
1	C	64	LEU
1	C	120	HIS
1	C	182	ARG
1	C	196	GLU
1	D	36	ASP
1	D	58	ARG
1	D	76	LEU
1	D	105	ARG
1	D	131	SER
1	D	168	LYS
1	D	171	LEU

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

Mol	Chain	Res	Type
1	A	155	HIS
1	A	158	ASN
1	B	37	HIS
1	B	195	GLN
1	C	135	HIS
1	D	135	HIS

### 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

### 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	188/218 (86%)	0.50	19 (10%) <b>12</b> <b>11</b>	21, 37, 88, 108	0
1	B	191/218 (87%)	0.56	26 (13%) <b>7</b> <b>6</b>	22, 39, 104, 121	0
1	C	183/218 (83%)	0.63	22 (12%) <b>9</b> <b>8</b>	19, 42, 96, 122	0
1	D	189/218 (86%)	0.80	28 (14%) <b>5</b> <b>5</b>	23, 43, 101, 125	0
All	All	751/872 (86%)	0.62	95 (12%) <b>8</b> <b>7</b>	19, 40, 99, 125	0

All (95) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	33	TYR	11.4
1	B	197	ALA	6.2
1	A	197	ALA	6.2
1	D	108	ILE	6.0
1	B	114	SER	5.8
1	C	197	ALA	5.2
1	B	109	ALA	5.1
1	D	109	ALA	5.0
1	A	109	ALA	4.9
1	D	197	ALA	4.9
1	D	67	ILE	4.4
1	D	64	LEU	4.3
1	B	36	ASP	4.2
1	D	107	TYR	4.1
1	B	80	CYS	3.9
1	B	64	LEU	3.9
1	D	169	VAL	3.6
1	D	115	GLY	3.6
1	C	118	SER	3.6
1	C	149	PHE	3.5
1	D	30	LEU	3.4

*Continued on next page...*

*Continued from previous page...*

<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	A	64	LEU	3.4
1	B	62	ILE	3.3
1	C	42	PHE	3.3
1	D	116	TYR	3.2
1	A	42	PHE	3.2
1	B	111	HIS	3.2
1	A	116	TYR	3.2
1	A	30	LEU	3.1
1	A	7	GLU	3.1
1	C	147	MET	3.0
1	D	62	ILE	3.0
1	B	113	GLU	2.9
1	D	57	ALA	2.9
1	B	115	GLY	2.9
1	C	62	ILE	2.8
1	D	5	LYS	2.8
1	B	107	TYR	2.8
1	B	6	TRP	2.8
1	A	107	TYR	2.7
1	C	159	TYR	2.7
1	B	4	ASP	2.7
1	C	4	ASP	2.7
1	A	115	GLY	2.7
1	D	156	SER	2.6
1	C	50	VAL	2.6
1	A	70	MET	2.6
1	D	65	HIS	2.6
1	C	64	LEU	2.6
1	D	3	ASP	2.6
1	C	58	ARG	2.6
1	C	34	GLU	2.5
1	B	105	ARG	2.5
1	D	171	LEU	2.5
1	D	173	LEU	2.5
1	C	36	ASP	2.4
1	B	32	GLU	2.4
1	B	46	ARG	2.4
1	C	54	LEU	2.4
1	D	68	GLU	2.4
1	B	108	ILE	2.4
1	B	117	ARG	2.4
1	D	59	ARG	2.4

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	D	63	PRO	2.4
1	A	51	ALA	2.4
1	B	30	LEU	2.3
1	D	152	THR	2.3
1	A	35	ASP	2.3
1	B	67	ILE	2.3
1	A	67	ILE	2.3
1	C	30	LEU	2.3
1	A	6	TRP	2.3
1	B	61	SER	2.2
1	B	52	SER	2.2
1	A	120	HIS	2.2
1	C	196	GLU	2.2
1	D	193	GLU	2.2
1	C	57	ALA	2.2
1	B	58	ARG	2.2
1	D	50	VAL	2.1
1	C	102	VAL	2.1
1	C	3	ASP	2.1
1	A	195	GLN	2.1
1	A	59	ARG	2.1
1	D	58	ARG	2.1
1	A	3	ASP	2.1
1	D	35	ASP	2.1
1	A	63	PRO	2.1
1	D	165	ILE	2.1
1	B	3	ASP	2.0
1	C	161	TYR	2.0
1	D	161	TYR	2.0
1	B	65	HIS	2.0
1	B	60	LYS	2.0
1	C	60	LYS	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

## 6.3 Carbohydrates [i](#)

There are no oligosaccharides 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.