



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 10:17 pm GMT

PDB ID : 3ZTH  
Title : Crystal structure of Stu0660 of Streptococcus thermophilus  
Authors : Lee, K.-H.; Robinson, H.; Oh, B.-H.  
Deposited on : 2011-07-08  
Resolution : 2.40 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  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
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) : recalc28949

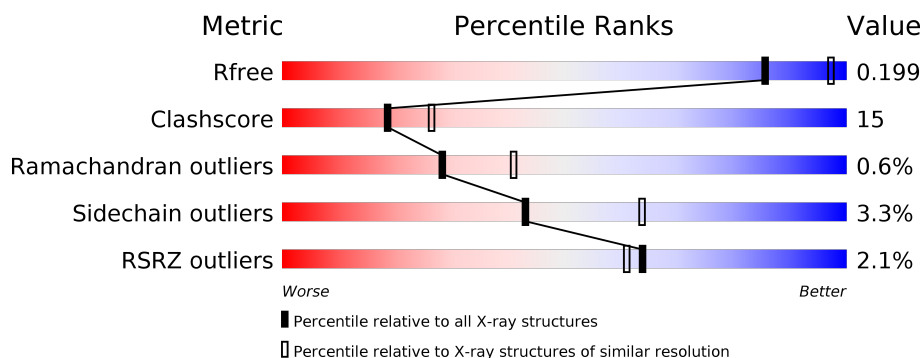
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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))
$R_{free}$	100719	3166 (2.40-2.40)
Clashscore	112137	3674 (2.40-2.40)
Ramachandran outliers	110173	3616 (2.40-2.40)
Sidechain outliers	110143	3617 (2.40-2.40)
RSRZ outliers	101464	3195 (2.40-2.40)

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	A	350	<div> <div>2%</div> <div> <div></div> <div>69%</div> <div>29%</div> <div>..</div> </div> </div>
1	B	350	<div> <div>2%</div> <div> <div></div> <div>70%</div> <div>27%</div> <div>..</div> </div> </div>

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
2	MG	A	1349	-	-	-	X
2	MG	B	1349	-	-	-	X

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6075 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 STU0660.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	348	Total	C	N	O	S	Se	0	0	1
			2917	1892	469	545	3	8			
1	B	348	Total	C	N	O	S	Se	0	0	1
			2917	1892	469	545	3	8			

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Mg	0	0
			1	1		
2	A	1	Total	Mg	0	0
			1	1		

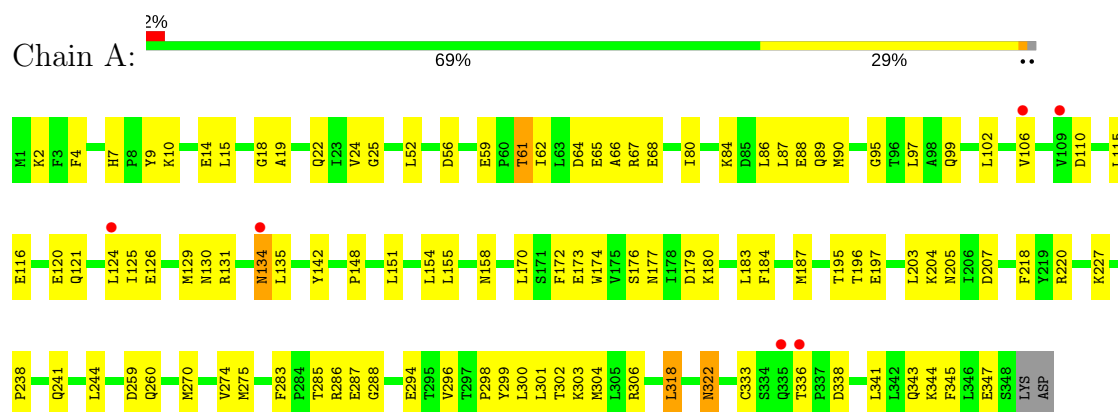
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	135	Total	O	0	0
			135	135		
3	B	104	Total	O	0	0
			104	104		

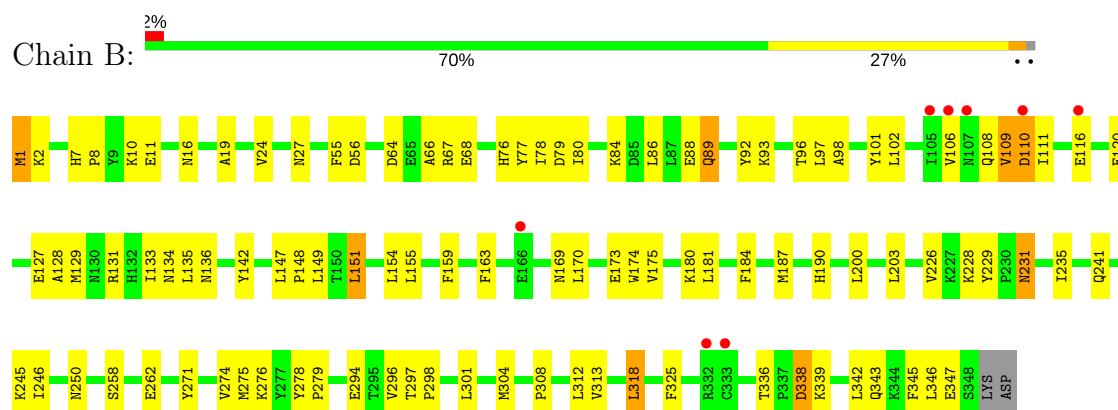
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: STU0660



#### • Molecule 1: STU0660



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	118.15Å 101.87Å 102.78Å 90.00° 115.14° 90.00°	Depositor
Resolution (Å)	50.00 – 2.40 45.88 – 2.40	Depositor EDS
% Data completeness (in resolution range)	98.9 (50.00-2.40) 98.1 (45.88-2.40)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.52 (at 2.39Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, $R_{free}$	0.205 , 0.243 0.203 , 0.199	Depositor DCC
$R_{free}$ test set	2144 reflections (5.31%)	DCC
Wilson B-factor (Å <sup>2</sup> )	37.3	Xtriage
Anisotropy	0.209	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 43.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	6075	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.65% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG

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	0.38	0/2978	0.61	0/4015
1	B	0.38	0/2978	0.62	1/4015 (0.0%)
All	All	0.38	0/5956	0.62	1/8030 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	76	HIS	N-CA-C	-5.04	97.40	111.00

There are no chirality outliers.

There are no planarity outliers.

### 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	2917	0	2891	97	0
1	B	2917	0	2891	83	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	135	0	0	7	0
3	B	104	0	0	2	0
All	All	6075	0	5782	172	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 15.

All (172) 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:285:THR:HG23	1:A:288:GLY:H	1.15	1.12
1:A:296:VAL:HG13	1:A:318:LEU:HD22	1.42	1.01
1:A:95:GLY:H	1:A:99:GLN:NE2	1.59	0.99
1:B:231:ASN:H	1:B:231:ASN:HD22	0.97	0.96
1:A:95:GLY:H	1:A:99:GLN:HE22	0.93	0.92
1:B:78:ILE:HD11	1:B:97:LEU:HD11	1.49	0.92
1:A:106:VAL:HG21	1:A:155:LEU:HD11	1.53	0.90
1:B:10:LYS:HG2	1:B:11:GLU:H	1.35	0.89
1:B:7:HIS:HD2	1:B:10:LYS:H	1.20	0.87
1:B:296:VAL:HG13	1:B:318:LEU:HD22	1.57	0.86
1:A:106:VAL:HG12	1:A:151:LEU:HD13	1.58	0.86
1:A:301:LEU:H	1:A:322:ASN:HD21	1.22	0.85
1:B:106:VAL:HG12	1:B:151:LEU:HD13	1.60	0.83
1:B:102:LEU:O	1:B:106:VAL:HG23	1.80	0.82
1:B:231:ASN:ND2	1:B:231:ASN:H	1.78	0.81
1:B:231:ASN:N	1:B:231:ASN:HD22	1.77	0.81
1:B:336:THR:HG22	1:B:338:ASP:OD1	1.79	0.81
1:B:19:ALA:HB1	1:B:226:VAL:HG11	1.63	0.81
1:A:97:LEU:HB3	1:A:187:MSE:HE3	1.63	0.80
1:A:195:THR:HG22	1:A:197:GLU:H	1.47	0.80
1:A:95:GLY:N	1:A:99:GLN:HE22	1.77	0.79
1:B:19:ALA:HB1	1:B:226:VAL:CG1	2.14	0.77
1:B:7:HIS:CD2	1:B:10:LYS:H	2.04	0.76
1:B:10:LYS:HG2	1:B:11:GLU:N	2.00	0.75
1:A:95:GLY:N	1:A:99:GLN:NE2	2.32	0.75
1:A:24:VAL:HG12	1:A:25:GLY:N	2.02	0.74
1:A:285:THR:HG22	3:A:2117:HOH:O	1.89	0.72
1:B:175:VAL:O	1:B:180:LYS:HE3	1.90	0.72
1:A:24:VAL:HG13	3:A:2018:HOH:O	1.89	0.72
1:B:1:MSE:HE2	1:B:2:LYS:H	1.54	0.72
1:A:343:GLN:HE22	1:B:339:LYS:NZ	1.88	0.71
1:A:285:THR:HG23	1:A:288:GLY:N	2.00	0.70
1:A:220:ARG:HD3	3:A:2099:HOH:O	1.92	0.69
1:B:250:ASN:HB2	3:B:2081:HOH:O	1.92	0.68
1:A:135:LEU:HD12	1:A:142:TYR:HD2	1.60	0.67
1:B:84:LYS:O	1:B:88:GLU:HG3	1.94	0.67
1:A:106:VAL:CG1	1:A:151:LEU:HD13	2.25	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:116:GLU:O	1:B:120:GLU:HG3	1.94	0.66
1:A:285:THR:CG2	1:A:288:GLY:H	2.02	0.65
1:A:116:GLU:O	1:A:120:GLU:HG3	1.97	0.64
1:B:7:HIS:HD2	1:B:10:LYS:N	1.95	0.64
1:A:345:PHE:CD2	1:B:279:PRO:HG3	2.33	0.64
3:A:2134:HOH:O	1:B:313:VAL:HG23	1.98	0.63
1:A:338:ASP:HB3	1:A:341:LEU:HD12	1.82	0.61
1:B:1:MSE:HE2	1:B:2:LYS:N	2.15	0.61
1:A:204:LYS:HG3	1:A:205:ASN:ND2	2.16	0.61
1:A:24:VAL:CG1	1:A:25:GLY:N	2.64	0.61
1:B:147:LEU:HD12	1:B:148:PRO:HD2	1.83	0.60
1:A:97:LEU:CB	1:A:187:MSE:HE3	2.34	0.58
1:B:86:LEU:HD22	1:B:184:PHE:CG	2.39	0.57
1:B:89:GLN:OE1	1:B:97:LEU:HD13	2.04	0.57
1:A:24:VAL:HG11	1:A:298:PRO:HG3	1.85	0.57
1:A:176:SER:HB3	1:A:179:ASP:OD2	2.05	0.57
1:A:66:ALA:O	1:A:67:ARG:HB2	2.04	0.56
1:A:84:LYS:HE2	1:A:88:GLU:OE2	2.05	0.56
1:B:77:TYR:CE2	1:B:79:ASP:HB2	2.41	0.56
1:B:271:TYR:O	1:B:275:MSE:HG2	2.06	0.55
1:B:149:LEU:HD21	1:B:154:LEU:HG	1.86	0.55
1:A:97:LEU:HD23	1:A:187:MSE:CE	2.36	0.55
1:A:125:ILE:O	1:A:129:MSE:HB2	2.07	0.54
1:B:262:GLU:HG2	1:B:325:PHE:CD1	2.43	0.54
1:B:19:ALA:CB	1:B:226:VAL:HG11	2.37	0.53
1:A:24:VAL:HG12	1:A:25:GLY:H	1.74	0.53
1:A:87:LEU:HB2	1:A:172:PHE:CE2	2.43	0.53
1:A:120:GLU:O	1:A:124:LEU:HG	2.07	0.53
1:A:300:LEU:HD23	1:A:318:LEU:HD13	1.91	0.53
1:A:106:VAL:HG11	1:A:151:LEU:HB3	1.90	0.53
1:A:22:GLN:OE1	1:A:244:LEU:HD12	2.09	0.52
1:A:135:LEU:HD12	1:A:142:TYR:CD2	2.43	0.52
1:A:97:LEU:HD23	1:A:187:MSE:HE3	1.90	0.52
1:A:7:HIS:HD2	1:A:10:LYS:H	1.57	0.52
1:B:155:LEU:HA	1:B:159:PHE:HD2	1.75	0.51
1:A:318:LEU:O	1:A:318:LEU:HD12	2.11	0.51
1:B:78:ILE:CD1	1:B:97:LEU:HD11	2.32	0.51
1:A:148:PRO:HG2	3:A:2067:HOH:O	2.11	0.50
1:B:127:GLU:O	1:B:131:ARG:HG3	2.10	0.50
1:A:343:GLN:HE22	1:B:339:LYS:CE	2.24	0.50
1:A:296:VAL:CG1	1:A:318:LEU:HD22	2.29	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:24:VAL:CG1	1:A:25:GLY:H	2.23	0.50
1:B:274:VAL:HG12	1:B:275:MSE:HE2	1.93	0.50
1:A:299:TYR:OH	1:B:241:GLN:NE2	2.45	0.50
1:B:10:LYS:CG	1:B:11:GLU:H	2.14	0.49
1:B:89:GLN:HG3	1:B:97:LEU:HD22	1.93	0.49
1:A:177:ASN:HA	1:A:180:LYS:HE2	1.94	0.49
1:B:108:GLN:O	1:B:110:ASP:N	2.46	0.49
1:A:121:GLN:O	1:A:125:ILE:HG13	2.12	0.49
1:B:101:TYR:CZ	1:B:190:HIS:HB3	2.48	0.49
1:A:336:THR:HG21	1:B:313:VAL:HG22	1.93	0.49
1:A:302:THR:HG22	1:A:304:MSE:HG3	1.96	0.48
1:A:154:LEU:HD23	1:A:158:ASN:HD22	1.78	0.48
1:A:322:ASN:HD22	1:A:322:ASN:HA	1.51	0.48
1:A:24:VAL:HG12	1:A:298:PRO:HA	1.96	0.48
1:A:97:LEU:CG	1:A:187:MSE:HE3	2.43	0.48
1:B:24:VAL:HG21	1:B:298:PRO:HG3	1.95	0.48
1:B:134:ASN:O	1:B:135:LEU:HD23	2.14	0.48
1:A:296:VAL:HG11	1:A:318:LEU:HA	1.95	0.47
1:B:10:LYS:HE2	1:B:11:GLU:HG2	1.96	0.47
1:A:196:THR:HG22	1:A:196:THR:O	2.13	0.47
1:A:195:THR:HG22	1:A:197:GLU:N	2.24	0.47
1:A:203:LEU:HD13	1:A:218:PHE:HZ	1.79	0.47
1:A:345:PHE:CG	1:B:279:PRO:HG3	2.49	0.47
1:A:344:LYS:O	1:A:347:GLU:HB2	2.15	0.47
1:B:339:LYS:O	1:B:343:GLN:HG3	2.14	0.47
1:B:109:VAL:O	1:B:111:ILE:N	2.47	0.47
1:A:318:LEU:C	1:A:318:LEU:HD12	2.35	0.46
1:A:2:LYS:HB2	1:A:65:GLU:HG3	1.97	0.46
1:B:129:MSE:O	1:B:133:ILE:HG13	2.16	0.46
1:B:19:ALA:HB1	1:B:226:VAL:HG13	1.92	0.46
1:B:342:LEU:O	1:B:345:PHE:HB3	2.15	0.46
1:A:18:GLY:O	1:A:19:ALA:C	2.54	0.46
1:B:245:LYS:HA	1:B:294:GLU:OE2	2.15	0.46
1:A:259:ASP:OD2	1:A:303:LYS:N	2.36	0.46
1:A:64:ASP:OD2	1:A:68:GLU:HB2	2.15	0.45
1:A:86:LEU:HD22	1:A:172:PHE:HZ	1.80	0.45
1:B:92:TYR:CD1	1:B:102:LEU:HD22	2.51	0.45
1:A:102:LEU:O	1:A:106:VAL:HG23	2.16	0.45
1:B:24:VAL:HG22	3:B:2018:HOH:O	2.15	0.45
1:A:227:LYS:HD2	3:A:2102:HOH:O	2.15	0.45
1:B:55:PHE:O	1:B:56:ASP:HB2	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:174:TRP:HA	1:A:174:TRP:CE3	2.51	0.45
1:A:64:ASP:HB2	1:A:65:GLU:OE1	2.17	0.44
1:A:207:ASP:OD2	1:A:238:PRO:HA	2.17	0.44
1:A:89:GLN:O	1:A:187:MSE:HE1	2.17	0.44
1:B:278:TYR:HA	1:B:279:PRO:HD3	1.79	0.44
1:B:98:ALA:HB2	1:B:187:MSE:HE3	2.00	0.44
1:B:262:GLU:HB2	1:B:301:LEU:HD21	2.00	0.44
1:B:173:GLU:N	1:B:173:GLU:OE1	2.49	0.44
1:B:297:THR:HB	1:B:298:PRO:HD3	1.99	0.44
1:A:287:GLU:HA	1:A:287:GLU:OE1	2.18	0.44
1:B:77:TYR:HE2	1:B:79:ASP:HB2	1.81	0.44
1:B:93:LYS:O	1:B:96:THR:HG23	2.18	0.44
1:B:163:PHE:O	1:B:169:ASN:HA	2.18	0.44
1:B:86:LEU:HD11	1:B:181:LEU:HD23	2.00	0.44
1:B:66:ALA:O	1:B:67:ARG:HB2	2.18	0.43
1:A:61:THR:HG23	1:A:62:ILE:N	2.32	0.43
1:A:333:CYS:HB2	1:B:346:LEU:HB2	2.01	0.43
1:B:246:ILE:HG12	1:B:294:GLU:HG2	2.00	0.43
1:B:296:VAL:HG13	1:B:318:LEU:CD2	2.40	0.43
1:B:64:ASP:OD1	1:B:66:ALA:N	2.49	0.43
1:B:128:ALA:HA	1:B:131:ARG:NH1	2.34	0.43
1:A:90:MSE:HG2	1:A:187:MSE:CE	2.49	0.43
1:A:52:LEU:HD22	1:A:59:GLU:HA	2.00	0.43
1:B:200:LEU:HD11	1:B:235:ILE:HG13	2.01	0.43
1:B:228:LYS:HE2	1:B:229:TYR:OH	2.18	0.43
1:A:270:MSE:O	1:A:274:VAL:HG23	2.19	0.43
1:B:27:ASN:HB2	1:B:258:SER:HA	2.01	0.43
1:A:296:VAL:HG11	1:A:318:LEU:HB2	2.00	0.42
1:B:80:ILE:HD12	1:B:86:LEU:HD23	2.01	0.42
1:A:126:GLU:HG2	1:A:130:ASN:ND2	2.35	0.42
1:A:4:PHE:CE1	1:A:14:GLU:HG3	2.54	0.42
1:A:90:MSE:HA	1:A:187:MSE:HE1	2.02	0.42
1:B:174:TRP:HA	1:B:174:TRP:CE3	2.55	0.42
1:A:173:GLU:N	1:A:173:GLU:OE1	2.46	0.42
1:B:135:LEU:HB2	1:B:142:TYR:HB3	2.01	0.42
1:A:87:LEU:HB2	1:A:172:PHE:CD2	2.54	0.42
1:A:90:MSE:HE1	1:A:183:LEU:HB3	2.02	0.42
1:A:244:LEU:O	1:A:294:GLU:HG2	2.20	0.41
1:A:174:TRP:HA	1:A:174:TRP:HE3	1.84	0.41
1:A:343:GLN:HG3	1:B:312:LEU:HD21	2.02	0.41
1:A:24:VAL:CG1	1:A:298:PRO:HA	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:ARG:O	1:A:134:ASN:OD1	2.38	0.41
1:B:7:HIS:ND1	1:B:8:PRO:HD2	2.35	0.41
1:A:220:ARG:HH11	1:A:220:ARG:HG2	1.85	0.41
1:A:296:VAL:HG11	1:A:318:LEU:CB	2.50	0.41
1:A:275:MSE:HG3	1:A:283:PHE:CD1	2.56	0.41
1:A:306:ARG:NH2	3:A:2123:HOH:O	2.43	0.41
1:B:276:LYS:HG2	1:B:276:LYS:O	2.21	0.40
1:B:102:LEU:HG	1:B:106:VAL:CG2	2.52	0.40
1:B:64:ASP:OD2	1:B:68:GLU:HB2	2.22	0.40
1:A:61:THR:CG2	1:A:62:ILE:N	2.84	0.40
1:B:304:MSE:HE1	1:B:308:PRO:HG3	2.03	0.40
1:A:80:ILE:HD11	1:A:184:PHE:CZ	2.56	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	346/350 (99%)	334 (96%)	11 (3%)	1 (0%)	44	60
1	B	346/350 (99%)	332 (96%)	11 (3%)	3 (1%)	20	29
All	All	692/700 (99%)	666 (96%)	22 (3%)	4 (1%)	28	41

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	109	VAL
1	B	110	ASP
1	A	134	ASN
1	B	347	GLU

### 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	331/326 (102%)	319 (96%)	12 (4%)	40	60
1	B	331/326 (102%)	321 (97%)	10 (3%)	46	67
All	All	662/652 (102%)	640 (97%)	22 (3%)	43	64

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	TYR
1	A	15	LEU
1	A	56	ASP
1	A	61	THR
1	A	110	ASP
1	A	115	LEU
1	A	170	LEU
1	A	241	GLN
1	A	260	GLN
1	A	286	ARG
1	A	318	LEU
1	A	322	ASN
1	B	1	MSE
1	B	16	ASN
1	B	89	GLN
1	B	136	ASN
1	B	151	LEU
1	B	170	LEU
1	B	203	LEU
1	B	231	ASN
1	B	318	LEU
1	B	338	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	7	HIS
1	A	76	HIS
1	A	99	GLN
1	A	134	ASN
1	A	143	HIS
1	A	153	GLN
1	A	158	ASN
1	A	205	ASN
1	A	241	GLN
1	A	281	ASN
1	A	322	ASN
1	A	323	GLN
1	A	343	GLN
1	B	7	HIS
1	B	16	ASN
1	B	153	GLN
1	B	231	ASN
1	B	241	GLN
1	B	260	GLN
1	B	263	HIS

### 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 carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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 [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, 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	340/350 (97%)	-0.26	6 (1%) 69 66	21, 42, 71, 84	0
1	B	340/350 (97%)	-0.16	8 (2%) 59 56	28, 44, 81, 98	0
All	All	680/700 (97%)	-0.21	14 (2%) 64 61	21, 43, 74, 98	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	106	VAL	5.3
1	A	335	GLN	4.7
1	A	109	VAL	3.9
1	A	134	ASN	3.1
1	A	336	THR	2.9
1	B	110	ASP	2.9
1	B	333	CYS	2.5
1	A	106	VAL	2.4
1	B	332	ARG	2.4
1	B	166	GLU	2.3
1	B	116	GLU	2.3
1	A	124	LEU	2.2
1	B	105	ILE	2.1
1	B	107	ASN	2.1

### 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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	MG	B	1349	1/1	0.96	0.28	6.86	35,35,35,35	0
2	MG	A	1349	1/1	0.98	0.23	4.79	38,38,38,38	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.