



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 10:27 PM GMT

PDB ID : 5D7R  
Title : Crystal structure of the ATP binding domain of *S. aureus* GyrB complexed with a ligand  
Authors : Zhang, J.; Yang, Q.; Cross, J.B.; Romero, J.A.C.; Ryan, M.D.; Lippa, B.; Dolle, R.E.; Andersen, O.A.; Barker, J.; Cheng, R.K.; Kahmann, J.; Felicetti, B.; Wood, M.; Scheich, C.  
Deposited on : 2015-08-14  
Resolution : 1.55 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 : 1.7.1 (RC1), CSD as537be (2016)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026982  
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) : rb-20026982

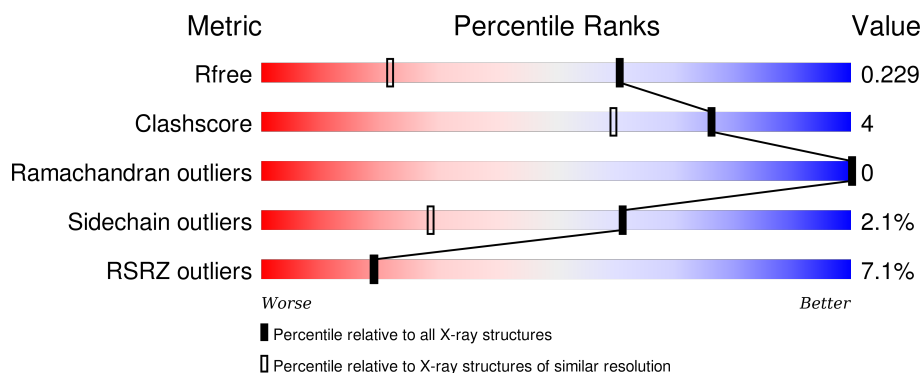
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.55 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	1665 (1.58-1.54)
Clashscore	102246	1014 (1.56-1.56)
Ramachandran outliers	100387	1704 (1.58-1.54)
Sidechain outliers	100360	1702 (1.58-1.54)
RSRZ outliers	91569	1668 (1.58-1.54)

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	212	
1	B	212	

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	B	304	-	-	-	X
4	57Y	A	306	-	-	-	X

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 3613 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 DNA gyrase subunit B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	193	Total	C	N	O	S	0	17	0
			1666	1048	291	325	2			
1	B	187	Total	C	N	O	S	0	10	0
			1564	987	269	306	2			

There are 50 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP P0A0K8
A	1	SER	-	expression tag	UNP P0A0K8
A	?	-	VAL	deletion	UNP P0A0K8
A	?	-	LEU	deletion	UNP P0A0K8
A	?	-	HIS	deletion	UNP P0A0K8
A	?	-	ALA	deletion	UNP P0A0K8
A	?	-	GLY	deletion	UNP P0A0K8
A	?	-	GLY	deletion	UNP P0A0K8
A	?	-	LYS	deletion	UNP P0A0K8
A	?	-	PHE	deletion	UNP P0A0K8
A	?	-	GLY	deletion	UNP P0A0K8
A	?	-	GLY	deletion	UNP P0A0K8
A	?	-	GLY	deletion	UNP P0A0K8
A	?	-	GLY	deletion	UNP P0A0K8
A	?	-	GLY	deletion	UNP P0A0K8
A	?	-	TYR	deletion	UNP P0A0K8
A	?	-	LYS	deletion	UNP P0A0K8
A	?	-	VAL	deletion	UNP P0A0K8
A	?	-	SER	deletion	UNP P0A0K8
A	?	-	GLY	deletion	UNP P0A0K8
A	?	-	GLY	deletion	UNP P0A0K8
A	?	-	LEU	deletion	UNP P0A0K8
A	?	-	HIS	deletion	UNP P0A0K8
A	?	-	GLY	deletion	UNP P0A0K8
A	?	-	VAL	deletion	UNP P0A0K8
A	?	-	GLY	deletion	UNP P0A0K8

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Chain	Residue	Modelled	Actual	Comment	Reference
B	0	GLY	-	expression tag	UNP P0A0K8
B	1	SER	-	expression tag	UNP P0A0K8
B	?	-	VAL	deletion	UNP P0A0K8
B	?	-	LEU	deletion	UNP P0A0K8
B	?	-	HIS	deletion	UNP P0A0K8
B	?	-	ALA	deletion	UNP P0A0K8
B	?	-	GLY	deletion	UNP P0A0K8
B	?	-	GLY	deletion	UNP P0A0K8
B	?	-	LYS	deletion	UNP P0A0K8
B	?	-	PHE	deletion	UNP P0A0K8
B	?	-	GLY	deletion	UNP P0A0K8
B	?	-	GLY	deletion	UNP P0A0K8
B	?	-	GLY	deletion	UNP P0A0K8
B	?	-	GLY	deletion	UNP P0A0K8
B	?	-	TYR	deletion	UNP P0A0K8
B	?	-	LYS	deletion	UNP P0A0K8
B	?	-	VAL	deletion	UNP P0A0K8
B	?	-	SER	deletion	UNP P0A0K8
B	?	-	GLY	deletion	UNP P0A0K8
B	?	-	GLY	deletion	UNP P0A0K8
B	?	-	LEU	deletion	UNP P0A0K8
B	?	-	HIS	deletion	UNP P0A0K8
B	?	-	GLY	deletion	UNP P0A0K8
B	?	-	VAL	deletion	UNP P0A0K8
B	?	-	GLY	deletion	UNP P0A0K8

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

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	4	Total Mg 4 4	0	0
2	A	4	Total Mg 4 4	0	0

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Cl 1 1	0	0
3	A	1	Total Cl 1 1	0	0

- 
- Chemical structure of compound 57Y, showing a complex heterocyclic system with multiple rings and substituents. The structure includes a central benzothiazine-like core, a 2-hydroxyethyl group (C1-C3), and a 2-hydroxy-3-(2-hydroxyethyl)acrylate group (C14-C18, O17, O18, O19). The atoms are color-coded: carbons in green, nitrogens in blue, sulfur in yellow, and oxygens in red.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 28	C 19	N 4	O 4	S 1	0	0
4	B	1	Total 28	C 19	N 4	O 4	S 1	0	0

- 
- Chemical structure of (S)-2-methylpentan-3-ol. The structure is labeled with green text: C1, C2, C3, C4(S), C5, O2, O4, and CM. The hydroxyl groups are shown in red as OH. The structure is a 3D representation with a central carbon (C2) bonded to a methyl group (CM), a hydroxyl group (OH, O2), and two other groups. One of these groups is a 1-hydroxyethyl group (C1, O4, H), and the other is a 1-methylpropyl group (C3, C4(S), C5).

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			8	6	2		

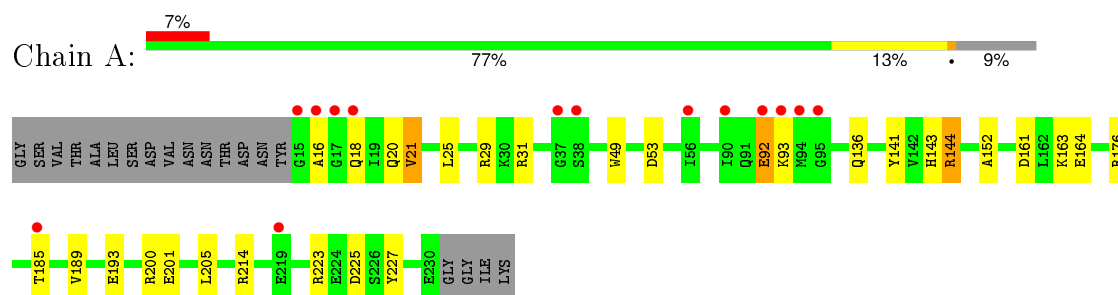
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	161	Total	O	0	0
			161	161		
6	B	148	Total	O	0	0
			148	148		

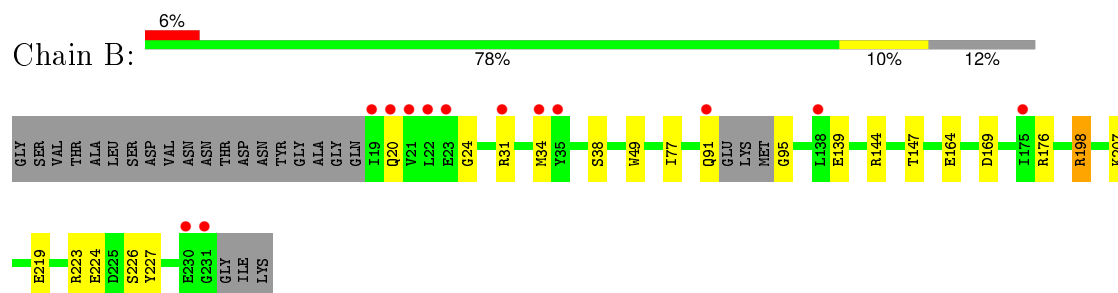
### 3 Residue-property plots [i](#)

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 ( $\text{RSRZ} > 2$ ). 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: DNA gyrase subunit B



#### • Molecule 1: DNA gyrase subunit B





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	143.19Å 55.51Å 51.11Å 90.00° 100.65° 90.00°	Depositor
Resolution (Å)	31.34 – 1.55 31.34 – 1.55	Depositor EDS
% Data completeness (in resolution range)	84.8 (31.34-1.55) 84.8 (31.34-1.55)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.12 (at 1.55Å)	Xtriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.181 , 0.220 0.193 , 0.229	Depositor DCC
$R_{free}$ test set	2445 reflections (5.30%)	DCC
Wilson B-factor (Å <sup>2</sup> )	25.0	Xtriage
Anisotropy	0.175	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 46.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 48606 reflections	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	3613	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.00% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 57Y, MPD, MG, CL

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	1.38	5/1737 (0.3%)	1.30	14/2346 (0.6%)
1	B	1.36	6/1619 (0.4%)	1.25	5/2189 (0.2%)
All	All	1.37	11/3356 (0.3%)	1.27	19/4535 (0.4%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	227	TYR	CE1-CZ	7.19	1.47	1.38
1	A	164	GLU	C-O	5.86	1.34	1.23
1	B	219	GLU	CD-OE1	5.58	1.31	1.25
1	A	143	HIS	N-CA	-5.48	1.35	1.46
1	B	139	GLU	CD-OE2	5.45	1.31	1.25
1	B	224	GLU	CD-OE2	-5.44	1.19	1.25
1	B	226	SER	CB-OG	-5.44	1.35	1.42
1	A	49	TRP	CB-CG	-5.13	1.41	1.50
1	A	163	LYS	CA-CB	-5.11	1.42	1.53
1	B	169	ASP	C-O	5.05	1.32	1.23
1	A	227	TYR	CE1-CZ	-5.02	1.32	1.38

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	169	ASP	CB-CG-OD1	11.88	128.99	118.30
1	A	176	ARG	NE-CZ-NH2	9.42	125.01	120.30
1	A	214	ARG	NE-CZ-NH2	-7.61	116.50	120.30
1	B	144	ARG	NE-CZ-NH2	-7.34	116.63	120.30
1	A	214	ARG	NE-CZ-NH1	7.09	123.85	120.30
1	A	144	ARG	NE-CZ-NH1	6.93	123.76	120.30
1	B	198	ARG	NE-CZ-NH1	6.28	123.44	120.30
1	B	223	ARG	NE-CZ-NH1	5.99	123.29	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	144	ARG	NE-CZ-NH2	-5.84	117.38	120.30
1	A	53	ASP	CB-CG-OD2	-5.81	113.07	118.30
1	A	152	ALA	N-CA-CB	5.54	117.85	110.10
1	A	141	TYR	CB-CG-CD2	-5.53	117.68	121.00
1	A	25	LEU	CB-CG-CD1	-5.41	101.81	111.00
1	A	141	TYR	CG-CD1-CE1	-5.39	116.99	121.30
1	A	189	VAL	CA-CB-CG2	-5.38	102.83	110.90
1	A	31	ARG	NE-CZ-NH1	-5.29	117.66	120.30
1	A	161	ASP	CB-CG-OD1	5.24	123.02	118.30
1	B	169	ASP	OD1-CG-OD2	-5.16	113.49	123.30
1	A	29	ARG	NE-CZ-NH1	5.07	122.83	120.30

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	1666	0	1672	18	0
1	B	1564	0	1558	12	0
2	A	4	0	0	0	0
2	B	4	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	28	0	14	2	0
4	B	28	0	15	0	0
5	B	8	0	14	0	0
6	A	161	0	0	7	0
6	B	148	0	0	2	0
All	All	3613	0	3273	25	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 4.

All (25) 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:223[C]:ARG:NH1	6:A:403:HOH:O	1.75	1.16
1:A:136[B]:GLN:OE1	6:A:402:HOH:O	1.62	1.16
1:A:185[B]:THR:O	6:A:404:HOH:O	1.76	1.02
1:A:20[B]:GLN:OE1	1:B:34[B]:MET:HE1	1.70	0.92
1:A:20[B]:GLN:OE1	1:B:34[B]:MET:CE	2.25	0.83
1:B:207:LYS:HG2	6:B:424:HOH:O	1.97	0.65
1:B:91:GLN:O	1:B:95:GLY:N	2.31	0.64
1:A:185[B]:THR:HB	6:A:404:HOH:O	2.02	0.59
1:A:18:GLN:HA	1:A:21:VAL:HG13	1.85	0.58
1:A:223[B]:ARG:HG3	6:A:407:HOH:O	2.02	0.57
1:A:200[A]:ARG:NH1	6:A:405:HOH:O	2.30	0.57
1:A:92:GLU:HG2	1:B:24:GLY:HA2	1.92	0.52
1:A:223[C]:ARG:HG3	6:A:407:HOH:O	2.12	0.48
1:B:207:LYS:CG	6:B:424:HOH:O	2.57	0.47
1:A:144:ARG:HH12	4:A:306:57Y:C18	2.27	0.45
1:B:49:TRP:CE3	1:B:49:TRP:HA	2.53	0.43
1:A:144:ARG:NH1	4:A:306:57Y:N	2.50	0.43
1:B:147:THR:HG23	1:B:164[B]:GLU:HG3	1.99	0.43
1:A:92:GLU:CG	1:B:24:GLY:HA2	2.48	0.43
1:A:223[B]:ARG:NH1	1:A:225:ASP:OD2	2.53	0.42
1:A:16:ALA:HB3	1:B:34[A]:MET:HG2	2.03	0.41
1:B:77:ILE:O	1:B:176:ARG:HA	2.20	0.41
1:B:147:THR:CG2	1:B:164[B]:GLU:HG3	2.51	0.41
1:A:201:GLU:O	1:A:205:LEU:HG	2.21	0.41
1:A:193:GLU:OE2	1:A:223[C]:ARG:NH2	2.53	0.41

There are no symmetry-related clashes.

## 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	209/212 (99%)	202 (97%)	7 (3%)	0	<b>100</b> <b>100</b>

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	193/212 (91%)	187 (97%)	6 (3%)	0	100	100
All	All	402/424 (95%)	389 (97%)	13 (3%)	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	186/183 (102%)	183 (98%)	3 (2%)	70	41
1	B	174/183 (95%)	169 (97%)	5 (3%)	50	17
All	All	360/366 (98%)	352 (98%)	8 (2%)	61	27

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	VAL
1	A	92	GLU
1	A	93	LYS
1	B	20	GLN
1	B	31	ARG
1	B	38[A]	SER
1	B	38[B]	SER
1	B	198	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	20	GLN
1	B	210	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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 13 ligands modelled in this entry, 10 are monoatomic - leaving 3 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z  > 2$	Counts	RMSZ	# $ Z  > 2$
4	57Y	A	306	-	23,31,31	1.43	5 (21%)	21,45,45	3.33	10 (47%)
5	MPD	B	305	-	6,7,7	0.63	0	6,10,10	1.30	1 (16%)
4	57Y	B	307	-	23,31,31	1.21	3 (13%)	21,45,45	3.26	14 (66%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	57Y	A	306	-	-	0/6/15/15	0/4/4/4
5	MPD	B	305	-	-	0/5/5/5	0/0/0/0
4	57Y	B	307	-	-	0/6/15/15	0/4/4/4

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	307	57Y	C16-CA	-2.93	1.36	1.40
4	A	306	57Y	C16-CA	-2.29	1.37	1.40
4	B	307	57Y	C20-C19	-2.24	1.37	1.42
4	A	306	57Y	C15-C14	-2.02	1.36	1.39
4	B	307	57Y	C10-C9	2.05	1.41	1.37
4	A	306	57Y	C10-N11	2.23	1.39	1.36
4	A	306	57Y	C3-C4	2.45	1.55	1.52
4	A	306	57Y	C22-C4	3.56	1.41	1.37

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	306	57Y	C22-C4-C5	-7.82	111.12	118.57
4	B	307	57Y	O17-C16-CA	-5.52	112.55	121.51
4	B	307	57Y	C22-C4-C5	-5.20	113.61	118.57
4	A	306	57Y	C2-C3-C4	-4.90	104.19	112.74
4	B	307	57Y	C21-C20-C19	-3.61	116.39	121.15
4	B	307	57Y	C18-C14-C12	-2.91	117.63	121.15
4	B	307	57Y	C14-C15-C16	-2.79	118.24	120.55
4	A	306	57Y	C21-C20-C19	-2.69	117.60	121.15
4	A	306	57Y	C15-C14-C18	-2.67	114.61	117.20
5	B	305	MPD	CM-C2-C1	-2.51	104.41	110.41
4	B	307	57Y	O23-C22-C4	-2.39	114.81	119.64
4	B	307	57Y	C18-N-CA	-2.37	114.33	118.22
4	A	306	57Y	C18-C14-C12	2.02	123.59	121.15
4	B	307	57Y	C10-N11-C12	2.05	108.93	103.98
4	B	307	57Y	C14-C18-N	2.17	127.82	124.31
4	A	306	57Y	C14-C18-N	2.18	127.85	124.31
4	B	307	57Y	C20-C19-C5	2.65	122.49	117.03
4	B	307	57Y	C3-C4-C22	2.97	125.78	120.25
4	A	306	57Y	C9-C10-N11	3.11	115.34	109.09
4	A	306	57Y	C10-N11-C12	3.39	112.16	103.98
4	B	307	57Y	O17-C16-C15	3.52	128.81	119.44
4	B	307	57Y	C9-C10-N11	4.30	117.72	109.09
4	A	306	57Y	C15-C16-CA	4.79	122.99	118.36
4	B	307	57Y	C21-C22-C4	7.88	125.78	121.44
4	A	306	57Y	C21-C22-C4	8.09	125.90	121.44

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	306	57Y	2	0

## 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	193/212 (91%)	0.56	14 (7%)	18 18	16, 22, 37, 72	0
1	B	187/212 (88%)	0.60	13 (6%)	19 19	15, 22, 40, 64	0
All	All	380/424 (89%)	0.58	27 (7%)	19 19	15, 22, 40, 72	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	19	ILE	7.4
1	A	16	ALA	6.9
1	B	230	GLU	6.9
1	B	21	VAL	5.9
1	A	92	GLU	5.0
1	B	35	TYR	5.0
1	A	94	MET	5.0
1	B	22	LEU	4.8
1	B	20	GLN	4.7
1	A	15	GLY	4.2
1	B	231	GLY	4.0
1	A	93	LYS	3.6
1	B	34[A]	MET	3.6
1	A	17	GLY	3.4
1	A	185[A]	THR	3.4
1	A	56	ILE	3.1
1	A	90	ILE	2.9
1	B	23	GLU	2.8
1	B	91	GLN	2.7
1	A	95	GLY	2.7
1	A	219	GLU	2.6
1	A	18	GLN	2.5
1	B	31	ARG	2.4
1	A	37	GLY	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	175[A]	ILE	2.3
1	A	38[A]	SER	2.2
1	B	138	LEU	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(Å <sup>2</sup> )	Q<0.9
2	MG	B	304	1/1	0.99	0.14	9.93	28,28,28,28	0
4	57Y	A	306	28/28	0.69	0.20	3.51	17,31,40,44	28
4	57Y	B	307	28/28	0.86	0.21	1.75	18,22,28,30	28
2	MG	A	304	1/1	0.95	0.16	1.40	44,44,44,44	0
5	MPD	B	305	8/8	0.79	0.14	1.00	29,35,38,40	0
2	MG	B	303	1/1	0.96	0.13	0.43	30,30,30,30	0
2	MG	A	303	1/1	0.97	0.03	-	45,45,45,45	1
2	MG	B	301	1/1	0.93	0.12	-	36,36,36,36	0
3	CL	A	305	1/1	0.98	0.03	-	38,38,38,38	0
2	MG	B	302	1/1	0.97	0.20	-	40,40,40,40	0
2	MG	A	302	1/1	0.96	0.13	-	30,30,30,30	0
3	CL	B	306	1/1	0.95	0.04	-	40,40,40,40	0
2	MG	A	301	1/1	0.77	0.19	-	67,67,67,67	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.