



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:28 PM GMT

PDB ID : 1XA7  
Title : Crystal structure of the benzylpenicillin-acylated BlaR1 sensor domain from *Staphylococcus aureus*  
Authors : Wilke, M.S.; Hills, T.L.; Zhang, H.Z.; Chambers, H.F.; Strynadka, N.C.  
Deposited on : 2004-08-25  
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  
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

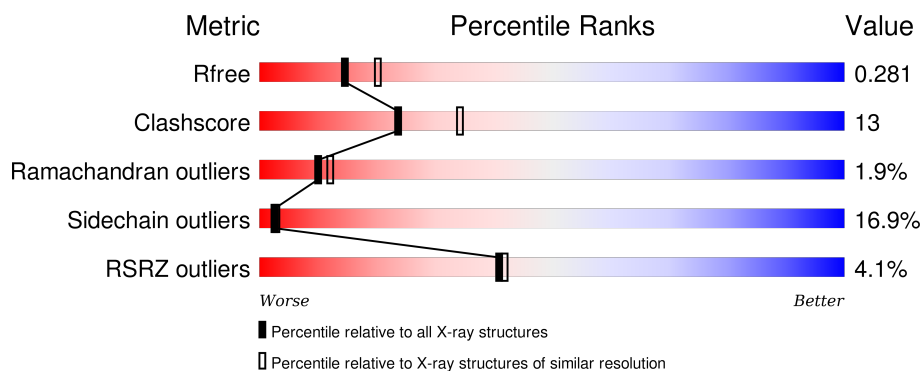
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*


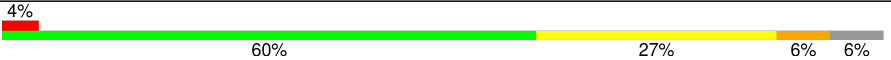
The reported resolution of this entry is 2.40 Å.

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))
$R_{free}$	91344	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (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	255	
1	B	255	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4106 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 Regulatory protein blaR1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	242	Total	C	N	O	S	Se	0	0	0
			2036	1304	341	381	1	9			
1	B	240	Total	C	N	O	S	Se	0	0	0
			2029	1301	341	377	1	9			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	35	MSE	MET	MODIFIED RESIDUE	UNP P18357
A	38	MSE	MET	MODIFIED RESIDUE	UNP P18357
A	59	PG1	SER	MODIFIED RESIDUE	UNP P18357
A	67	MSE	MET	MODIFIED RESIDUE	UNP P18357
A	82	MSE	MET	MODIFIED RESIDUE	UNP P18357
A	104	MSE	MET	MODIFIED RESIDUE	UNP P18357
A	146	MSE	MET	MODIFIED RESIDUE	UNP P18357
A	164	MSE	MET	MODIFIED RESIDUE	UNP P18357
A	165	MSE	MET	MODIFIED RESIDUE	UNP P18357
A	249	MSE	MET	MODIFIED RESIDUE	UNP P18357
B	35	MSE	MET	MODIFIED RESIDUE	UNP P18357
B	38	MSE	MET	MODIFIED RESIDUE	UNP P18357
B	59	PG1	SER	MODIFIED RESIDUE	UNP P18357
B	67	MSE	MET	MODIFIED RESIDUE	UNP P18357
B	82	MSE	MET	MODIFIED RESIDUE	UNP P18357
B	104	MSE	MET	MODIFIED RESIDUE	UNP P18357
B	146	MSE	MET	MODIFIED RESIDUE	UNP P18357
B	164	MSE	MET	MODIFIED RESIDUE	UNP P18357
B	165	MSE	MET	MODIFIED RESIDUE	UNP P18357
B	249	MSE	MET	MODIFIED RESIDUE	UNP P18357

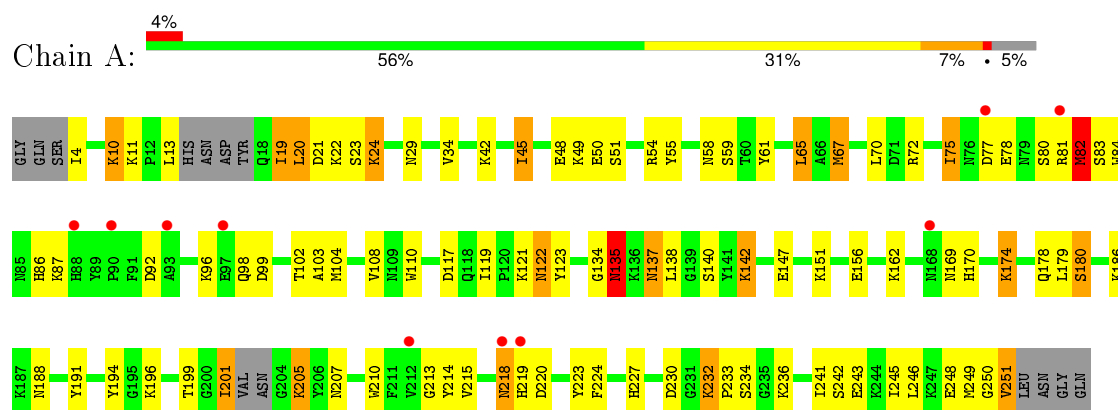
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	21	Total 21	O 21	0	0
2	B	20	Total 20	O 20	0	0

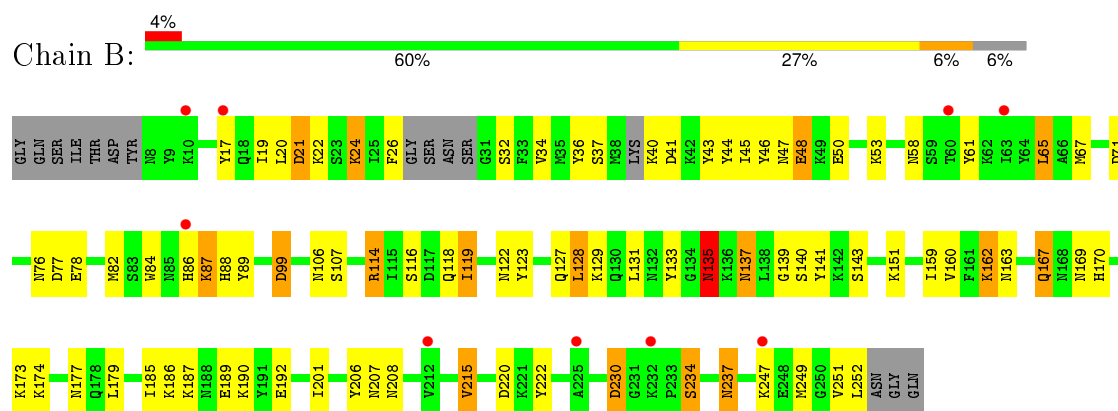
### 3 Residue-property plots

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: Regulatory protein blaR1



#### • Molecule 1: Regulatory protein blaR1



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	88.39 Å 88.39 Å 125.06 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.40 30.32 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.1 (30.00-2.40) 99.1 (30.32-2.00)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.87 (at 2.00 Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, $R_{free}$	0.220 , 0.275 0.224 , 0.281	Depositor DCC
$R_{free}$ test set	971 reflections (5.14%)	DCC
Wilson B-factor (Å <sup>2</sup> )	34.8	Xtriage
Anisotropy	0.158	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 46.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.43$ , $\langle L^2 \rangle = 0.26$	Xtriage
Outliers	0 of 33840 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	4106	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 3.45% 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

Bond lengths and bond angles in the following residue types are not validated in this section: PG1

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	0.92	3/2045 (0.1%)	0.96	6/2727 (0.2%)
1	B	0.87	1/2039 (0.0%)	0.92	8/2721 (0.3%)
All	All	0.89	4/4084 (0.1%)	0.94	14/5448 (0.3%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	249	MSE	CG-SE	-5.85	1.75	1.95
1	A	223	TYR	CD2-CE2	-5.44	1.31	1.39
1	A	210	TRP	CB-CG	-5.37	1.40	1.50
1	A	67	MSE	SE-CE	-5.03	1.65	1.95

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	99	ASP	CB-CG-OD2	8.42	125.88	118.30
1	B	220	ASP	CB-CG-OD2	7.86	125.38	118.30
1	A	21	ASP	CB-CG-OD2	7.25	124.83	118.30
1	A	230	ASP	CB-CG-OD2	7.19	124.77	118.30
1	A	220	ASP	CB-CG-OD2	6.56	124.20	118.30
1	B	71	ASP	CB-CG-OD2	6.53	124.18	118.30
1	B	21	ASP	CB-CG-OD2	5.93	123.64	118.30
1	B	230	ASP	CB-CG-OD2	5.89	123.60	118.30
1	B	99	ASP	CB-CG-OD2	5.82	123.54	118.30
1	B	77	ASP	CB-CG-OD2	5.45	123.21	118.30
1	A	92	ASP	CB-CG-OD2	5.44	123.20	118.30
1	B	41	ASP	CB-CG-OD2	5.44	123.20	118.30
1	A	117	ASP	CB-CG-OD2	5.31	123.08	118.30
1	B	114	ARG	NE-CZ-NH1	5.02	122.81	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	2036	0	1985	60	0
1	B	2029	0	1974	46	0
2	A	21	0	0	3	0
2	B	20	0	0	3	0
All	All	4106	0	3959	106	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 13.

All (106) 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:98:GLN:HE22	1:A:102:THR:HG22	1.29	0.96
1:A:98:GLN:HE22	1:A:102:THR:CG2	1.80	0.95
1:A:67:MSE:HE3	1:A:179:LEU:HD22	1.50	0.93
1:A:82:MSE:SE	2:A:271:HOH:O	2.36	0.92
1:A:98:GLN:NE2	1:A:102:THR:HG22	1.90	0.87
1:A:98:GLN:NE2	1:A:102:THR:CG2	2.41	0.84
1:A:174:LYS:HD3	1:A:178:GLN:HE21	1.46	0.81
1:B:190:LYS:NZ	1:B:252:LEU:O	2.15	0.76
1:B:22:LYS:HA	1:B:24:LYS:HZ2	1.49	0.75
1:A:242:SER:O	1:A:246:LEU:HD23	1.91	0.70
1:A:55:TYR:CD2	1:A:227:HIS:CE1	2.82	0.68
1:B:187:LYS:NZ	1:B:192:GLU:OE2	2.25	0.67
1:A:180:SER:OG	1:A:214:TYR:OH	2.12	0.67
1:A:137:ASN:C	1:A:137:ASN:HD22	1.97	0.67
1:A:29:ASN:HD22	1:A:232:LYS:H	1.43	0.67
1:B:118:GLN:HG3	2:B:270:HOH:O	1.94	0.66
1:A:45:ILE:HD11	1:A:48:GLU:HA	1.79	0.65
1:A:137:ASN:ND2	1:A:138:LEU:O	2.31	0.63
1:A:67:MSE:CE	1:A:179:LEU:HD22	2.28	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:98:GLN:HG2	1:A:110:TRP:CZ2	2.34	0.63
1:B:24:LYS:HD3	1:B:24:LYS:H	1.64	0.62
1:A:218:ASN:HD22	1:A:218:ASN:H	1.48	0.62
1:A:135:ASN:HB3	1:A:137:ASN:H	1.64	0.62
1:A:98:GLN:HE22	1:A:102:THR:HG21	1.64	0.62
1:A:201:ILE:HA	1:A:205:LYS:O	2.00	0.61
1:B:67:MSE:HE2	1:B:179:LEU:HD22	1.84	0.60
1:B:22:LYS:HA	1:B:24:LYS:NZ	2.16	0.60
1:B:84:TRP:CZ2	1:B:86:HIS:HA	2.38	0.59
1:B:36:TYR:CE2	1:B:162:LYS:HD2	2.38	0.59
1:B:37:SER:OG	1:B:40:LYS:HB2	2.02	0.59
1:A:29:ASN:ND2	1:A:232:LYS:H	2.01	0.58
1:B:67:MSE:CE	1:B:179:LEU:HD22	2.32	0.58
1:B:34:VAL:HG22	1:B:45:ILE:HG23	1.85	0.57
1:A:70:LEU:HD23	1:A:75:ILE:HG12	1.87	0.57
1:B:123:TYR:O	1:B:127:GLN:HG2	2.05	0.56
1:B:67:MSE:CE	1:B:179:LEU:CD2	2.83	0.56
1:A:84:TRP:CE2	1:A:86:HIS:HA	2.41	0.56
1:A:20:LEU:HD21	1:A:249:MSE:CE	2.36	0.55
1:A:55:TYR:CE2	1:A:227:HIS:CE1	2.94	0.55
1:B:163:ASN:OD1	1:B:167:GLN:OE1	2.25	0.54
1:A:188:ASN:ND2	1:A:191:TYR:CE1	2.72	0.54
1:A:142:LYS:CE	1:A:142:LYS:HA	2.38	0.53
1:B:118:GLN:CG	2:B:270:HOH:O	2.56	0.53
1:B:207:ASN:ND2	1:B:208:ASN:HB2	2.24	0.53
1:A:98:GLN:NE2	1:A:102:THR:HG21	2.20	0.53
1:A:174:LYS:HD3	1:A:178:GLN:NE2	2.19	0.53
1:A:19:ILE:HB	1:A:45:ILE:O	2.09	0.53
1:B:185:ILE:HG22	1:B:186:LYS:HG3	1.91	0.53
1:B:159:ILE:HD13	1:B:159:ILE:N	2.24	0.53
1:A:10:LYS:H	1:A:10:LYS:CD	2.23	0.52
1:A:34:VAL:O	1:A:224:PHE:HA	2.10	0.52
1:A:54:ARG:HB3	1:A:151:LYS:HB3	1.92	0.51
1:A:10:LYS:HD2	1:A:10:LYS:H	1.75	0.51
1:A:241:ILE:O	1:A:245:ILE:HG12	2.10	0.50
1:A:134:GLY:HA2	1:A:156:GLU:OE2	2.12	0.50
1:A:194:TYR:O	1:A:213:GLY:HA3	2.12	0.50
1:B:46:TYR:O	1:B:47:ASN:HB2	2.12	0.49
1:B:106:ASN:ND2	2:B:263:HOH:O	2.25	0.49
1:A:98:GLN:HG2	1:A:110:TRP:CH2	2.47	0.49
1:B:67:MSE:HE3	1:B:179:LEU:CD2	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:103:ALA:HA	1:A:108:VAL:HG23	1.95	0.48
1:A:59:PG1:H12	1:A:199:THR:HB	1.96	0.48
1:A:137:ASN:ND2	1:A:137:ASN:C	2.67	0.47
1:B:45:ILE:HG21	1:B:48:GLU:HB2	1.95	0.47
1:B:61:TYR:CD2	1:B:65:LEU:HD22	2.50	0.47
1:B:43:TYR:HB3	1:B:45:ILE:HD11	1.95	0.47
1:B:234:SER:HB3	1:B:237:ASN:H	1.80	0.46
1:A:207:ASN:HA	1:A:233:PRO:HG3	1.96	0.46
1:A:22:LYS:HA	1:A:24:LYS:HE3	1.96	0.46
1:A:67:MSE:HE1	1:A:179:LEU:HB2	1.98	0.46
1:B:206:TYR:CD2	1:B:206:TYR:N	2.84	0.45
1:B:137:ASN:HD22	1:B:139:GLY:H	1.64	0.45
1:A:232:LYS:HD2	1:A:232:LYS:HA	1.87	0.45
1:A:142:LYS:NZ	1:A:142:LYS:HA	2.31	0.45
1:B:67:MSE:HE3	1:B:179:LEU:HD23	1.98	0.45
1:B:163:ASN:OD1	1:B:167:GLN:CD	2.56	0.44
1:A:45:ILE:HD11	1:A:48:GLU:CA	2.46	0.44
1:A:121:LYS:NZ	2:A:258:HOH:O	2.49	0.44
1:A:250:GLY:O	1:A:251:VAL:O	2.35	0.44
1:A:61:TYR:CD2	1:A:65:LEU:HD22	2.53	0.44
1:A:77:ASP:HB2	1:A:174:LYS:HZ3	1.84	0.43
1:B:140:SER:O	1:B:141:TYR:HB3	2.17	0.43
1:B:135:ASN:HB2	1:B:151:LYS:H	1.84	0.43
1:B:116:SER:HA	1:B:119:ILE:HG13	1.99	0.43
1:A:186:LYS:NZ	1:A:243:GLU:OE2	2.51	0.43
1:B:114:ARG:O	1:B:118:GLN:NE2	2.52	0.43
1:A:142:LYS:HA	1:A:142:LYS:HE2	1.99	0.43
1:B:206:TYR:N	1:B:206:TYR:HD2	2.17	0.43
1:A:82:MSE:CG	2:A:271:HOH:O	2.67	0.43
1:B:87:LYS:O	1:B:89:TYR:N	2.52	0.42
1:B:20:LEU:HD11	1:B:44:TYR:HB3	2.02	0.42
1:B:177:ASN:HD22	1:B:177:ASN:N	2.18	0.42
1:B:131:LEU:O	1:B:160:VAL:HG13	2.20	0.42
1:B:128:LEU:HG	1:B:133:TYR:HB3	2.02	0.42
1:B:21:ASP:O	1:B:24:LYS:HE3	2.20	0.41
1:B:207:ASN:HD22	1:B:208:ASN:HB2	1.85	0.41
1:B:215:VAL:HG13	1:B:222:TYR:HB2	2.02	0.41
1:A:98:GLN:CG	1:A:110:TRP:CZ2	3.02	0.41
1:A:104:MSE:SE	1:A:196:LYS:HD3	2.70	0.41
1:B:82:MSE:HB3	1:B:114:ARG:NH2	2.35	0.41
1:B:135:ASN:HB3	1:B:137:ASN:H	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:122:ASN:HD22	1:A:123:TYR:N	2.19	0.41
1:A:72:ARG:HD3	1:A:119:ILE:HG12	2.01	0.41
1:A:135:ASN:OD1	1:A:137:ASN:HB3	2.20	0.40
1:A:45:ILE:HD11	1:A:48:GLU:N	2.37	0.40
1:B:251:VAL:HG12	1:B:251:VAL:O	2.21	0.40

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	235/255 (92%)	210 (89%)	19 (8%)	6 (3%)	7	6
1	B	233/255 (91%)	213 (91%)	17 (7%)	3 (1%)	15	21
All	All	468/510 (92%)	423 (90%)	36 (8%)	9 (2%)	10	12

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	135	ASN
1	A	23	SER
1	B	58	ASN
1	A	51	SER
1	A	58	ASN
1	A	82	MSE
1	B	88	HIS
1	A	147	GLU
1	B	135	ASN

### 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	220/222 (99%)	180 (82%)	40 (18%)	2	2
1	B	219/222 (99%)	185 (84%)	34 (16%)	3	4
All	All	439/444 (99%)	365 (83%)	74 (17%)	2	3

All (74) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	ILE
1	A	10	LYS
1	A	11	LYS
1	A	13	LEU
1	A	19	ILE
1	A	20	LEU
1	A	24	LYS
1	A	42	LYS
1	A	45	ILE
1	A	49	LYS
1	A	50	GLU
1	A	65	LEU
1	A	75	ILE
1	A	78	GLU
1	A	80	SER
1	A	81	ARG
1	A	82	MSE
1	A	83	SER
1	A	87	LYS
1	A	96	LYS
1	A	122	ASN
1	A	135	ASN
1	A	137	ASN
1	A	140	SER
1	A	142	LYS
1	A	162	LYS
1	A	169	ASN

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Mol	Chain	Res	Type
1	A	170	HIS
1	A	174	LYS
1	A	180	SER
1	A	201	ILE
1	A	205	LYS
1	A	215	VAL
1	A	218	ASN
1	A	219	HIS
1	A	232	LYS
1	A	234	SER
1	A	236	LYS
1	A	248	GLU
1	A	251	VAL
1	B	17	TYR
1	B	19	ILE
1	B	24	LYS
1	B	26	PHE
1	B	32	SER
1	B	48	GLU
1	B	50	GLU
1	B	53	LYS
1	B	65	LEU
1	B	76	ASN
1	B	78	GLU
1	B	87	LYS
1	B	99	ASP
1	B	107	SER
1	B	119	ILE
1	B	122	ASN
1	B	128	LEU
1	B	129	LYS
1	B	135	ASN
1	B	137	ASN
1	B	143	SER
1	B	162	LYS
1	B	167	GLN
1	B	169	ASN
1	B	170	HIS
1	B	173	LYS
1	B	174	LYS
1	B	189	GLU
1	B	201	ILE

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Mol	Chain	Res	Type
1	B	215	VAL
1	B	230	ASP
1	B	234	SER
1	B	237	ASN
1	B	247	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	29	ASN
1	A	98	GLN
1	A	122	ASN
1	A	137	ASN
1	A	154	ASN
1	A	167	GLN
1	A	178	GLN
1	A	218	ASN
1	B	76	ASN
1	B	105	GLN
1	B	135	ASN
1	B	137	ASN
1	B	177	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues are modelled in this entry.

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
1	PG1	A	59	1	23,30,31	1.89	4 (17%)	28,42,44	1.91	7 (25%)
1	PG1	B	59	1	23,30,31	1.98	5 (21%)	28,42,44	1.87	8 (28%)

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
1	PG1	A	59	1	-	0/17/42/44	0/2/2/2
1	PG1	B	59	1	-	0/17/42/44	0/2/2/2

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	59	PG1	C3-N4	-3.89	1.42	1.46
1	A	59	PG1	C2-S1	-3.77	1.77	1.85
1	B	59	PG1	C2-S1	-3.63	1.78	1.85
1	B	59	PG1	C3-N4	-2.18	1.44	1.46
1	B	59	PG1	C2-C3	-2.09	1.52	1.56
1	A	59	PG1	O-C	3.97	1.38	1.19
1	A	59	PG1	OG-C7	4.93	1.43	1.33
1	B	59	PG1	O-C	5.17	1.43	1.19
1	B	59	PG1	OG-C7	5.82	1.45	1.33

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	59	PG1	O-C-CA	-4.73	113.16	125.49
1	B	59	PG1	O-C-CA	-4.47	113.85	125.49
1	A	59	PG1	OG-C7-O8	-3.18	117.56	124.05
1	B	59	PG1	OG-C7-O8	-2.35	119.25	124.05
1	B	59	PG1	C9-C2-C3	-2.23	106.81	112.12
1	A	59	PG1	O16-C15-N14	-2.20	119.28	123.01
1	B	59	PG1	C18-C17-C15	-2.03	106.23	112.84
1	B	59	PG1	C17-C15-N14	2.01	119.19	115.96
1	A	59	PG1	C5-C6-N14	2.12	114.52	109.57
1	B	59	PG1	C10-C2-S1	2.14	112.75	109.17
1	A	59	PG1	C17-C15-N14	2.19	119.47	115.96
1	A	59	PG1	C2-S1-C5	2.20	98.71	93.98
1	B	59	PG1	C2-S1-C5	2.87	100.15	93.98
1	B	59	PG1	OG-C7-C6	5.49	118.83	111.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	59	PG1	OG-C7-C6	5.82	119.29	111.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	59	PG1	1	0

## 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 [i](#)

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	232/255 (90%)	0.12	10 (4%) 39 40	22, 44, 67, 85	0
1	B	230/255 (90%)	0.22	9 (3%) 43 44	24, 47, 71, 81	0
All	All	462/510 (90%)	0.17	19 (4%) 41 42	22, 46, 69, 85	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	77	ASP	5.0
1	A	88	HIS	3.3
1	A	219	HIS	3.2
1	A	218	ASN	3.2
1	B	232	LYS	3.2
1	B	60	THR	3.0
1	A	90	PRO	2.9
1	A	212	VAL	2.9
1	A	81	ARG	2.8
1	B	17	TYR	2.7
1	B	225	ALA	2.5
1	B	247	LYS	2.4
1	B	212	VAL	2.4
1	A	168	ASN	2.4
1	B	86	HIS	2.4
1	B	10	LYS	2.1
1	A	93	ALA	2.1
1	A	97	GLU	2.1
1	B	63	ILE	2.1

### 6.2 Non-standard residues in protein, DNA, RNA chains [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
1	PG1	A	59	29/30	0.91	0.15	-	23,34,46,48	0
1	PG1	B	59	29/30	0.93	0.14	-	23,32,42,45	0

### 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.