



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 10:33 PM GMT

PDB ID : 5DF8  
Title : CRYSTAL STRUCTURE OF PENICILLIN-BINDING PROTEIN 3 FROM  
PSEUDOMONAS AERUGINOSA IN COMPLEX WITH CEFOPERAZONE  
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Deposited on : 2015-08-26  
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  
<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

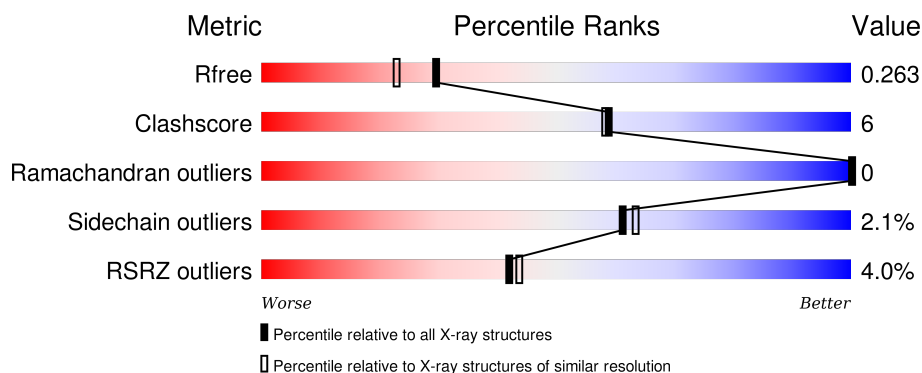
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 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))
$R_{free}$	91344	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (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 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	564	 2% 79% 11% • 9%
1	B	564	 5% 76% 10% • 9%

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
3	GOL	A	603	-	-	-	X
3	GOL	B	603	-	-	-	X

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8353 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 Cell division protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	514	Total	C	N	O	S	0	0	0
			3922	2474	712	724	12			
1	B	489	Total	C	N	O	S	0	0	0
			3717	2349	666	690	12			

There are 38 discrepancies between the modelled and reference sequences:

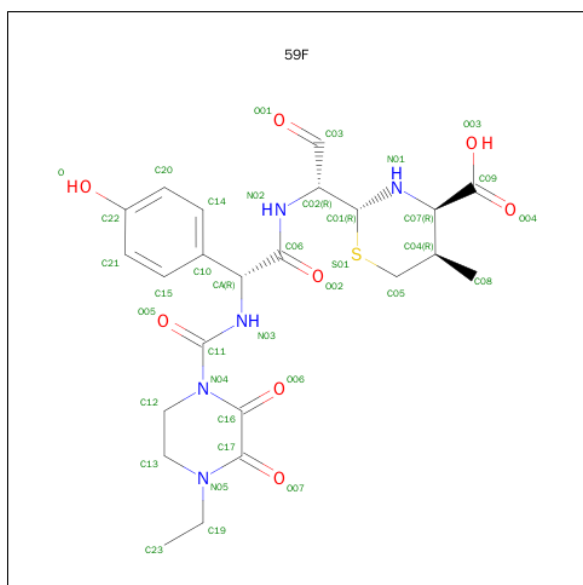
Chain	Residue	Modelled	Actual	Comment	Reference
A	16	MET	-	initiating methionine	UNP Q51504
A	17	ALA	-	expression tag	UNP Q51504
A	18	HIS	-	expression tag	UNP Q51504
A	19	HIS	-	expression tag	UNP Q51504
A	20	HIS	-	expression tag	UNP Q51504
A	21	HIS	-	expression tag	UNP Q51504
A	22	HIS	-	expression tag	UNP Q51504
A	23	HIS	-	expression tag	UNP Q51504
A	24	SER	-	expression tag	UNP Q51504
A	25	SER	-	expression tag	UNP Q51504
A	26	GLY	-	expression tag	UNP Q51504
A	27	LEU	-	expression tag	UNP Q51504
A	28	GLU	-	expression tag	UNP Q51504
A	29	VAL	-	expression tag	UNP Q51504
A	30	LEU	-	expression tag	UNP Q51504
A	31	PHE	-	expression tag	UNP Q51504
A	32	GLN	-	expression tag	UNP Q51504
A	33	GLY	-	expression tag	UNP Q51504
A	34	PRO	-	expression tag	UNP Q51504
B	16	MET	-	initiating methionine	UNP Q51504
B	17	ALA	-	expression tag	UNP Q51504
B	18	HIS	-	expression tag	UNP Q51504
B	19	HIS	-	expression tag	UNP Q51504
B	20	HIS	-	expression tag	UNP Q51504
B	21	HIS	-	expression tag	UNP Q51504

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Chain	Residue	Modelled	Actual	Comment	Reference
B	22	HIS	-	expression tag	UNP Q51504
B	23	HIS	-	expression tag	UNP Q51504
B	24	SER	-	expression tag	UNP Q51504
B	25	SER	-	expression tag	UNP Q51504
B	26	GLY	-	expression tag	UNP Q51504
B	27	LEU	-	expression tag	UNP Q51504
B	28	GLU	-	expression tag	UNP Q51504
B	29	VAL	-	expression tag	UNP Q51504
B	30	LEU	-	expression tag	UNP Q51504
B	31	PHE	-	expression tag	UNP Q51504
B	32	GLN	-	expression tag	UNP Q51504
B	33	GLY	-	expression tag	UNP Q51504
B	34	PRO	-	expression tag	UNP Q51504

- Molecule 2 is (2R,4R,5R)-2-[(1R)-1-[(2R)-2-[(4-ethyl-2,3-dioxopiperazin-1-yl)carbonyl]amino]-2-(4-hydroxyphenyl)acetyl]amino]-2-oxoethyl]-5-methyl-1,3-thiazinane-4-carboxylic acid (three-letter code: 59F) (formula: C<sub>23</sub>H<sub>29</sub>N<sub>5</sub>O<sub>8</sub>S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			37	23	5	8	1		
2	B	1	Total	C	N	O	S	0	0
			37	23	5	8	1		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 6 3 3	0	0
3	A	1	Total C O 6 3 3	0	0
3	A	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	316	Total O 316 316	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	282	Total 282	O 282	0	0





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	57.22Å 74.37Å 82.45Å 71.69° 86.06° 85.88°	Depositor
Resolution (Å)	47.21 – 2.00 47.17 – 2.00	Depositor EDS
% Data completeness (in resolution range)	92.8 (47.21-2.00) 84.1 (47.17-2.00)	Depositor EDS
$R_{merge}$	0.19	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.85 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, $R_{free}$	0.226 , 0.263 0.226 , 0.263	Depositor DCC
$R_{free}$ test set	4067 reflections (5.31%)	DCC
Wilson B-factor (Å <sup>2</sup> )	18.1	Xtriage
Anisotropy	0.330	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 57.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 80608 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	8353	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.80% 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: GOL, CL, 59F

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	0.37	0/3999	0.55	0/5426
1	B	0.37	0/3792	0.54	0/5147
All	All	0.37	0/7791	0.55	0/10573

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	A	3922	0	3983	46	0
1	B	3717	0	3755	44	0
2	A	37	0	25	1	0
2	B	37	0	25	1	0
3	A	18	0	24	2	0
3	B	18	0	24	1	0
4	A	3	0	0	0	0
4	B	3	0	0	0	0
5	A	316	0	0	5	0
5	B	282	0	0	8	0
All	All	8353	0	7836	90	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 (90) 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:478:GLY:H	1:A:562:THR:HG23	1.31	0.95
1:B:478:GLY:H	1:B:562:THR:HG23	1.32	0.90
1:B:395:ARG:HH11	1:B:395:ARG:HG3	1.37	0.87
1:B:319:TYR:HB2	1:B:338:ARG:HG3	1.61	0.82
1:A:319:TYR:HB2	1:A:338:ARG:HG3	1.64	0.79
1:A:286:MET:HG2	1:A:416:ILE:HD13	1.68	0.74
1:A:81:THR:HG22	1:A:127:ARG:HA	1.71	0.72
1:B:81:THR:HG22	1:B:128:GLY:H	1.54	0.72
1:B:432:VAL:HG13	1:B:443:PRO:HG2	1.71	0.72
1:A:395:ARG:HG2	1:A:395:ARG:HH11	1.56	0.70
1:A:432:VAL:HG13	1:A:443:PRO:HG2	1.74	0.70
1:B:81:THR:HG22	1:B:127:ARG:HA	1.75	0.69
1:A:189:VAL:HG12	1:A:214:LYS:HB2	1.75	0.69
1:A:173:ARG:NH1	5:A:702:HOH:O	2.26	0.68
1:B:81:THR:HG21	1:B:150:GLU:OE2	1.94	0.67
1:B:364:GLU:HG2	5:B:723:HOH:O	1.96	0.66
1:A:152:ARG:HH11	1:A:152:ARG:HG2	1.61	0.64
1:B:211:LYS:N	5:B:702:HOH:O	2.31	0.64
1:A:186:LEU:O	1:A:216:GLY:HA3	1.99	0.63
1:B:324:GLN:HG2	1:B:329:THR:HG22	1.81	0.61
1:B:562:THR:HG21	5:B:813:HOH:O	2.00	0.60
1:A:81:THR:HG21	1:A:150:GLU:OE2	2.06	0.55
1:B:286:MET:HG2	1:B:416:ILE:HD13	1.89	0.55
1:A:81:THR:CG2	1:A:150:GLU:OE2	2.56	0.54
1:B:81:THR:HG22	1:B:128:GLY:N	2.23	0.53
1:B:395:ARG:HH11	1:B:395:ARG:CG	2.17	0.52
1:A:560:LEU:N	1:A:561:PRO:HD2	2.24	0.52
1:B:395:ARG:HG3	1:B:395:ARG:NH1	2.15	0.52
1:B:250:VAL:HG21	1:B:416:ILE:HD12	1.91	0.52
1:B:222:SER:HB2	1:B:258:GLU:HB3	1.91	0.52
1:B:540:PRO:HB2	3:B:603:GOL:H31	1.92	0.52
1:B:81:THR:CG2	1:B:150:GLU:OE2	2.58	0.51
1:A:474:ALA:HB2	1:A:539:ALA:HB1	1.92	0.51
1:B:460:MET:O	1:B:464:VAL:HG23	2.10	0.51
1:B:86:PRO:O	1:B:90:MET:HG2	2.11	0.51
1:B:354:ILE:HG21	1:B:407:TYR:HB3	1.93	0.51
1:A:81:THR:HG22	1:A:128:GLY:H	1.77	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:80:THR:OG1	1:A:131:PRO:HA	2.11	0.50
1:B:498:TYR:CE1	2:B:601:59F:H2	2.46	0.50
1:A:301:MET:HG2	1:A:460:MET:HE1	1.93	0.49
1:A:284:ARG:HD2	5:A:750:HOH:O	2.11	0.49
1:A:499:ARG:HD2	5:A:730:HOH:O	2.12	0.49
1:A:96:TRP:HB2	1:A:97:PRO:HD3	1.94	0.49
1:A:395:ARG:NH1	1:A:395:ARG:HG2	2.25	0.48
1:A:282:ARG:HG2	3:A:603:GOL:H12	1.95	0.48
1:B:135:GLU:HA	1:B:138:ILE:HB	1.95	0.48
1:B:560:LEU:N	1:B:561:PRO:HD2	2.28	0.48
1:B:81:THR:HG22	1:B:127:ARG:CA	2.42	0.48
1:A:239:LEU:HD13	1:A:265:GLN:HB2	1.96	0.47
1:B:80:THR:OG1	1:B:131:PRO:HA	2.14	0.47
1:A:385:GLY:HA2	1:B:241:GLU:O	2.15	0.47
1:B:312:PRO:HB3	5:B:785:HOH:O	2.14	0.47
1:B:167:PHE:CE2	1:B:176:GLU:HG3	2.50	0.47
1:A:207:VAL:HG23	1:A:208:GLN:HG2	1.97	0.47
1:B:474:ALA:HB2	1:B:539:ALA:HB1	1.96	0.46
1:B:438:ARG:NH1	5:B:701:HOH:O	2.31	0.46
1:A:84:ALA:HA	1:A:144:GLY:O	2.16	0.46
1:A:56:ILE:HD11	1:A:193:ARG:CZ	2.45	0.45
1:A:56:ILE:HD11	1:A:193:ARG:NH1	2.31	0.45
1:A:106:ASP:OD2	1:A:109:LEU:HG	2.16	0.45
1:A:81:THR:HG21	1:A:127:ARG:HE	1.81	0.44
1:B:375:LEU:HD22	1:B:413:VAL:HG11	1.99	0.44
1:B:277:GLN:NE2	5:B:710:HOH:O	2.47	0.44
1:A:539:ALA:N	1:A:540:PRO:HD2	2.31	0.44
1:B:161:VAL:O	1:B:161:VAL:HG22	2.18	0.44
1:A:460:MET:O	1:A:464:VAL:HG23	2.16	0.44
1:B:324:GLN:CG	1:B:329:THR:HG22	2.48	0.43
1:A:139:ALA:O	1:A:141:LYS:HG2	2.18	0.43
1:A:274:ARG:HG2	1:A:275:ASN:OD1	2.18	0.43
1:A:83:TRP:CE2	1:A:146:TYR:HB2	2.52	0.43
1:A:432:VAL:CG1	1:A:443:PRO:HG2	2.45	0.43
1:A:265:GLN:HB3	5:A:732:HOH:O	2.18	0.43
1:A:479:TYR:CE1	1:A:557:PRO:HA	2.53	0.43
1:A:81:THR:HG22	1:A:127:ARG:CA	2.43	0.42
1:B:426:ALA:HB2	1:B:510:PHE:CG	2.54	0.42
1:A:233:ARG:HG2	5:A:950:HOH:O	2.19	0.42
1:A:375:LEU:HD22	1:A:413:VAL:HG11	2.00	0.42
1:A:282:ARG:CG	3:A:603:GOL:H12	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:539:ALA:N	1:B:540:PRO:HD2	2.34	0.41
1:B:96:TRP:HB2	1:B:97:PRO:HD3	2.02	0.41
1:B:432:VAL:CG1	1:B:443:PRO:HG2	2.45	0.41
1:B:245:LYS:HE3	5:B:859:HOH:O	2.19	0.41
1:A:378:ASP:HA	1:A:388:VAL:HG22	2.01	0.41
1:B:478:GLY:N	1:B:562:THR:HG23	2.16	0.41
2:A:601:59F:O06	2:A:601:59F:N03	2.44	0.41
1:A:492:SER:HA	1:A:499:ARG:HG2	2.02	0.41
1:B:348:LYS:HD3	5:B:808:HOH:O	2.21	0.41
1:B:479:TYR:CE1	1:B:557:PRO:HA	2.55	0.41
1:A:324:GLN:HG2	1:A:329:THR:HG22	2.03	0.41
1:A:71:GLU:HA	1:A:72:PRO:HD3	1.94	0.41

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	512/564 (91%)	500 (98%)	12 (2%)	0	100	100
1	B	485/564 (86%)	475 (98%)	10 (2%)	0	100	100
All	All	997/1128 (88%)	975 (98%)	22 (2%)	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	410/448 (92%)	401 (98%)	9 (2%)	60	62
1	B	387/448 (86%)	379 (98%)	8 (2%)	61	63
All	All	797/896 (89%)	780 (98%)	17 (2%)	61	63

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	61	HIS
1	A	67	ASP
1	A	80	THR
1	A	141	LYS
1	A	195	VAL
1	A	239	LEU
1	A	354	ILE
1	A	395	ARG
1	A	416	ILE
1	B	59	PRO
1	B	61	HIS
1	B	80	THR
1	B	103	LEU
1	B	149	GLU
1	B	354	ILE
1	B	393	ASN
1	B	395	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	277	GLN
1	B	393	ASN
1	B	475	GLN

### 5.3.3 RNA ⓘ

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 14 ligands modelled in this entry, 6 are monoatomic - leaving 8 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$
2	59F	A	601	1	31,39,39	1.95	8 (25%)	34,55,55	1.88	10 (29%)
3	GOL	A	602	-	5,5,5	0.25	0	5,5,5	0.57	0
3	GOL	A	603	-	5,5,5	0.32	0	5,5,5	0.44	0
3	GOL	A	604	-	5,5,5	0.26	0	5,5,5	0.21	0
2	59F	B	601	1	31,39,39	1.94	6 (19%)	34,55,55	1.79	6 (17%)
3	GOL	B	602	-	5,5,5	0.25	0	5,5,5	0.38	0
3	GOL	B	603	-	5,5,5	0.23	0	5,5,5	0.39	0
3	GOL	B	604	-	5,5,5	0.25	0	5,5,5	0.52	0

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
2	59F	A	601	1	-	0/21/62/62	0/2/3/3
3	GOL	A	602	-	-	0/4/4/4	0/0/0/0
3	GOL	A	603	-	-	0/4/4/4	0/0/0/0
3	GOL	A	604	-	-	0/4/4/4	0/0/0/0
2	59F	B	601	1	-	0/21/62/62	0/2/3/3
3	GOL	B	602	-	-	0/4/4/4	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	B	603	-	-	0/4/4/4	0/0/0/0
3	GOL	B	604	-	-	0/4/4/4	0/0/0/0

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	59F	C17-C16	-4.12	1.45	1.53
2	B	601	59F	C17-C16	-3.75	1.46	1.53
2	A	601	59F	C02-N02	-2.87	1.43	1.46
2	A	601	59F	C07-N01	-2.53	1.44	1.47
2	B	601	59F	C02-N02	-2.29	1.44	1.46
2	A	601	59F	CA-C06	-2.16	1.48	1.54
2	A	601	59F	C11-N04	2.78	1.47	1.42
2	B	601	59F	C11-N04	3.12	1.48	1.42
2	A	601	59F	C11-N03	3.97	1.44	1.34
2	A	601	59F	C17-N05	4.34	1.40	1.35
2	B	601	59F	C11-N03	4.35	1.45	1.34
2	B	601	59F	C17-N05	4.52	1.40	1.35
2	A	601	59F	C06-N02	4.61	1.44	1.34
2	B	601	59F	C06-N02	5.22	1.46	1.34

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	59F	O07-C17-N05	-2.60	120.48	123.56
2	A	601	59F	C15-C10-CA	-2.28	116.99	120.80
2	B	601	59F	O07-C17-N05	-2.22	120.93	123.56
2	A	601	59F	O01-C03-C02	-2.18	119.05	125.81
2	A	601	59F	C12-C13-N05	2.37	115.43	110.44
2	B	601	59F	C01-C02-N02	2.43	115.71	109.78
2	A	601	59F	O05-C11-N04	2.48	122.75	119.92
2	A	601	59F	C13-C12-N04	2.57	115.39	110.25
2	B	601	59F	C09-C07-N01	2.61	114.39	110.53
2	A	601	59F	C10-CA-C06	2.69	113.83	107.83
2	A	601	59F	C01-C02-N02	2.97	117.03	109.78
2	A	601	59F	C16-C17-N05	3.20	119.96	117.42
2	B	601	59F	C10-CA-C06	3.62	115.91	107.83
2	B	601	59F	N03-C11-N04	3.94	120.16	114.06
2	A	601	59F	C08-C04-C05	5.84	118.25	109.05
2	B	601	59F	C08-C04-C05	6.29	118.96	109.05

There are no chirality outliers.



There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	59F	1	0
3	A	603	GOL	2	0
2	B	601	59F	1	0
3	B	603	GOL	1	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	514/564 (91%)	-0.04	9 (1%) 71 72	8, 24, 60, 92	0
1	B	489/564 (86%)	0.15	31 (6%) 23 24	10, 26, 68, 102	0
All	All	1003/1128 (88%)	0.05	40 (3%) 42 44	8, 25, 65, 102	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	139	ALA	5.0
1	B	135	GLU	4.2
1	B	146	TYR	4.2
1	B	441	ARG	3.9
1	A	199	ARG	3.6
1	B	89	LEU	3.5
1	B	141	LYS	3.5
1	B	134	GLY	3.1
1	B	192	LYS	3.1
1	B	213	ALA	3.1
1	B	85	ASN	3.1
1	A	442	VAL	3.0
1	A	200	ARG	3.0
1	A	564	THR	2.9
1	B	148	ILE	2.9
1	B	57	ALA	2.9
1	B	327	ARG	2.8
1	B	119	GLU	2.8
1	A	440	ASP	2.7
1	A	441	ARG	2.7
1	B	143	PRO	2.7
1	B	102	ALA	2.7
1	B	136	GLY	2.7
1	A	561	PRO	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	87	LYS	2.5
1	B	80	THR	2.5
1	B	138	ILE	2.5
1	B	212	ASN	2.5
1	B	193	ARG	2.4
1	B	122	PHE	2.4
1	B	395	ARG	2.4
1	B	83	TRP	2.3
1	B	120	ARG	2.3
1	B	118	ALA	2.2
1	B	132	GLU	2.1
1	B	84	ALA	2.1
1	A	319	TYR	2.0
1	A	139	ALA	2.0
1	B	58	ILE	2.0
1	B	95	ARG	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 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
3	GOL	A	603	6/6	0.87	0.20	4.46	39,55,61,62	0
3	GOL	B	603	6/6	0.81	0.21	3.96	40,46,52,58	0
3	GOL	B	604	6/6	0.95	0.13	1.87	23,30,35,39	0
3	GOL	A	604	6/6	0.89	0.13	0.48	28,50,52,53	0
3	GOL	A	602	6/6	0.95	0.14	-0.13	13,19,27,27	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	59F	B	601	37/37	0.94	0.12	-0.29	12,23,33,38	0
3	GOL	B	602	6/6	0.95	0.13	-0.61	22,26,27,32	0
2	59F	A	601	37/37	0.96	0.09	-1.04	11,19,32,39	0
4	CL	B	605	1/1	0.98	0.06	-	38,38,38,38	0
4	CL	A	607	1/1	0.99	0.08	-	33,33,33,33	0
4	CL	B	607	1/1	0.95	0.05	-	51,51,51,51	0
4	CL	A	605	1/1	0.95	0.11	-	34,34,34,34	0
4	CL	A	606	1/1	0.98	0.05	-	48,48,48,48	0
4	CL	B	606	1/1	0.97	0.09	-	31,31,31,31	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.