



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 10:33 PM GMT

PDB ID : 5DF7
Title : CRYSTAL STRUCTURE OF PENICILLIN-BINDING PROTEIN 3 FROM
PSEUDOMONAS AERUGINOSA IN COMPLEX WITH AZLOCILLIN
Authors : Ren, J.; Nettleship, J.E.; Males, A.; Stuart, D.I.; Owens, R.J.
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
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

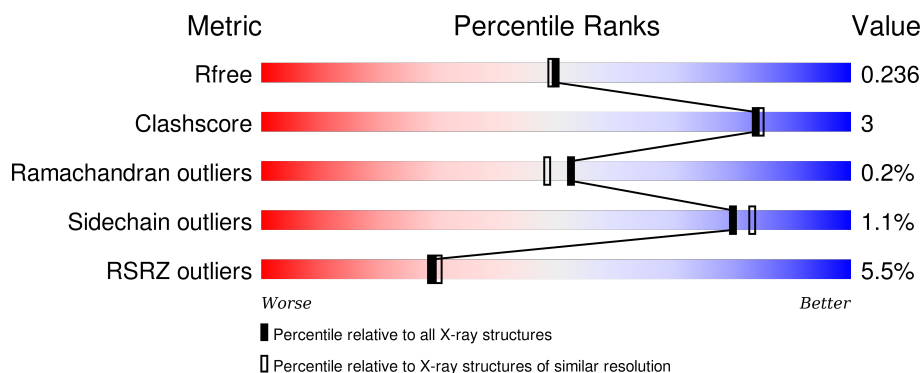
1 Overall quality at a glance

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 ≥ 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	
1	B	564	

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	B	602	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 8127 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	515	Total	C	N	O	S	0	0	0
			3927	2477	713	725	12			
1	B	490	Total	C	N	O	S	0	0	0
			3722	2352	667	691	12			

There are 38 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	16	MET	-	engineered mutation	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	-	engineered mutation	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

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- Chemical structure diagram showing a complex molecule with various labeled atoms and groups. The structure includes a benzene ring (labeled CX3, CX4, CX5, CX6, CX1, CX2) and a pyrazole ring (labeled CX8, CX9, CXA, CXB, CXE). Key functional groups include a carboxylic acid (O1, C1, O3, C3), an amide (N1, C, O), a sulfonamide (S4, O, CXC), and a pyrazole (N1, N3, CX1, CX2, CX3, CX4, CX5, CX6, CX7, CX8, CX9, CXA, CXB, CXE). The molecule is also labeled with 59H.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 32	C 20	N 5	O 6	S 1	0	0
2	B	1	Total 32	C 20	N 5	O 6	S 1	0	0

- WORLDWIDE
PDB
PROTEIN DATA BANK

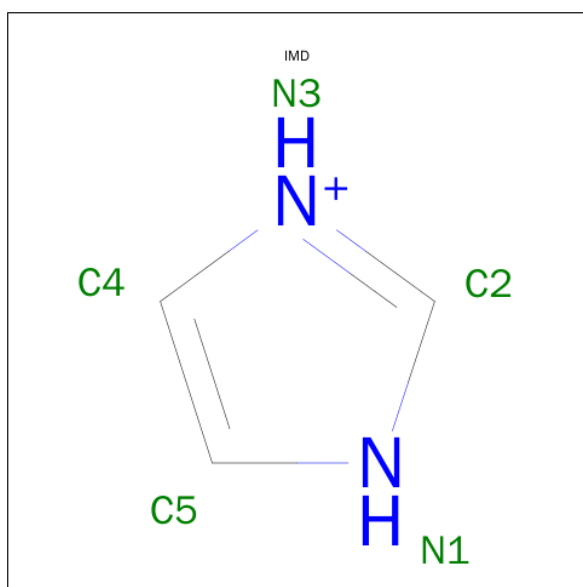


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		

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

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

- Molecule 5 is IMIDAZOLE (three-letter code: IMD) (formula: C₃H₅N₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	N	0	0
			5	3	2		
5	B	1	Total	C	N	0	0
			5	3	2		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	218	Total	O	0	0
			218	218		
6	B	164	Total	O	0	0
			164	164		

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	57.25Å 74.92Å 82.72Å 71.26° 85.99° 85.69°	Depositor
Resolution (Å)	47.19 – 2.00 47.19 – 2.00	Depositor EDS
% Data completeness (in resolution range)	86.7 (47.19-2.00) 78.9 (47.19-2.00)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.89 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.198 , 0.237 0.197 , 0.236	Depositor DCC
R_{free} test set	3803 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	26.2	Xtriage
Anisotropy	0.287	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 54.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 75952 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8127	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.21% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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, IMD, 59H

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	0.34	0/4004	0.54	0/5433
1	B	0.33	0/3797	0.52	0/5154
All	All	0.34	0/7801	0.53	0/10587

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	3927	0	3988	21	0
1	B	3722	0	3760	26	0
2	A	32	0	23	0	0
2	B	32	0	23	0	0
3	A	6	0	8	2	0
3	B	12	0	16	5	0
4	A	1	0	0	0	0
4	B	3	0	0	0	0
5	A	5	0	5	0	0
5	B	5	0	5	0	0
6	A	218	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	164	0	0	1	0
All	All	8127	0	7828	46	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 3.

All (46) 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:B:269:ASN:H	3:B:603:GOL:H31	1.56	0.70
1:B:321:GLY:HA3	1:B:335:ARG:HD3	1.78	0.66
1:B:562:THR:HG21	6:B:803:HOH:O	1.99	0.62
1:A:51:ARG:HH11	1:A:198:ASP:HB2	1.65	0.62
1:A:478:GLY:H	1:A:562:THR:HG23	1.66	0.60
1:A:518:ILE:HD12	1:A:550:LEU:HD23	1.84	0.59
1:B:478:GLY:H	1:B:562:THR:HG23	1.67	0.59
1:A:262:MET:SD	1:A:416:ILE:HD11	2.44	0.57
1:A:169:ASP:HB3	1:A:175:ARG:HG3	1.88	0.56
1:B:268:TYR:HB2	3:B:603:GOL:H32	1.88	0.55
1:B:322:THR:HG22	1:B:331:ARG:HG2	1.87	0.55
1:A:168:THR:HB	1:A:172:ASP:HA	1.88	0.54
1:A:301:MET:HG2	1:A:460:MET:HE1	1.90	0.53
1:B:86:PRO:O	1:B:90:MET:HG2	2.10	0.52
1:B:259:ILE:HD11	1:B:416:ILE:HD11	1.91	0.51
1:B:191:GLY:HA2	1:B:211:LYS:HG2	1.92	0.51
1:B:354:ILE:HG21	1:B:407:TYR:HB3	1.93	0.51
1:A:120:ARG:NH1	1:B:338:ARG:HD3	2.26	0.50
1:B:560:LEU:N	1:B:561:PRO:HD2	2.27	0.49
1:A:181:ALA:HB2	1:A:384:PRO:HD3	1.95	0.48
1:B:181:ALA:HB2	1:B:384:PRO:HD3	1.96	0.48
1:A:525:ASP:OD1	3:A:602:GOL:H11	2.14	0.47
1:B:173:ARG:HH11	1:B:175:ARG:HH11	1.62	0.47
1:B:523:VAL:HG11	3:B:602:GOL:H11	1.97	0.46
1:A:354:ILE:HG21	1:A:407:TYR:HB3	1.98	0.46
1:A:525:ASP:CG	3:A:602:GOL:H11	2.35	0.46
1:B:58:ILE:HD11	1:B:193:ARG:HB3	1.97	0.46
1:B:83:TRP:CE2	1:B:146:TYR:HB2	2.51	0.45
1:B:343:THR:HA	1:B:460:MET:HE2	1.98	0.45
1:A:51:ARG:NH1	1:A:198:ASP:HB2	2.31	0.44
1:A:86:PRO:O	1:A:90:MET:HG2	2.17	0.44
1:A:518:ILE:HD11	1:A:553:MET:SD	2.57	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:342:LEU:HA	1:B:345:ILE:HD12	2.00	0.44
1:B:256:THR:HG22	1:B:442:VAL:HG22	1.99	0.44
1:B:248:SER:HB2	3:B:602:GOL:H12	1.99	0.43
1:B:67:ASP:HB2	1:B:223:ILE:O	2.18	0.43
1:A:186:LEU:O	1:A:216:GLY:HA3	2.19	0.43
1:A:95:ARG:NH1	1:A:140:LEU:O	2.43	0.43
1:A:283:ASN:ND2	1:A:286:MET:HG3	2.34	0.42
1:A:222:SER:HB2	1:A:258:GLU:HB3	2.02	0.42
1:A:560:LEU:N	1:A:561:PRO:HD2	2.35	0.41
1:B:298:PRO:HG3	1:B:461:LEU:HD21	2.01	0.41
1:B:86:PRO:HG2	1:B:120:ARG:O	2.20	0.41
1:B:442:VAL:HA	1:B:443:PRO:HD3	1.88	0.40
1:B:269:ASN:H	3:B:603:GOL:C3	2.28	0.40
1:A:88:GLU:CD	1:A:202:ARG:HB3	2.41	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	513/564 (91%)	500 (98%)	12 (2%)	1 (0%)	52	48
1	B	486/564 (86%)	474 (98%)	11 (2%)	1 (0%)	52	48
All	All	999/1128 (89%)	974 (98%)	23 (2%)	2 (0%)	52	48

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	172	ASP
1	B	93	LYS

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%)	402 (98%)	8 (2%)	63	65
1	B	387/448 (86%)	386 (100%)	1 (0%)	94	96
All	All	797/896 (89%)	788 (99%)	9 (1%)	80	83

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	67	ASP
1	A	161	VAL
1	A	171	ASP
1	A	173	ARG
1	A	349	SER
1	A	354	ILE
1	A	416	ILE
1	A	441	ARG
1	B	175	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	212	ASN

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 [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 11 ligands modelled in this entry, 4 are monoatomic - leaving 7 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	59H	A	601	1	27,34,34	2.79	9 (33%)	34,49,49	1.83	8 (23%)
3	GOL	A	602	-	5,5,5	0.30	0	5,5,5	0.86	0
5	IMD	A	604	-	3,5,5	0.39	0	4,5,5	0.60	0
2	59H	B	601	1	27,34,34	2.85	9 (33%)	34,49,49	1.86	7 (20%)
3	GOL	B	602	-	5,5,5	0.16	0	5,5,5	0.53	0
3	GOL	B	603	-	5,5,5	0.26	0	5,5,5	0.36	0
5	IMD	B	607	-	3,5,5	0.38	0	4,5,5	0.62	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	59H	A	601	1	-	0/20/55/55	0/3/3/3
3	GOL	A	602	-	-	0/4/4/4	0/0/0/0
5	IMD	A	604	-	-	0/0/0/0	0/1/1/1
2	59H	B	601	1	-	0/20/55/55	0/3/3/3
3	GOL	B	602	-	-	0/4/4/4	0/0/0/0
3	GOL	B	603	-	-	0/4/4/4	0/0/0/0
5	IMD	B	607	-	-	0/0/0/0	0/1/1/1

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	601	59H	CXC-S4	-3.39	1.78	1.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	59H	CXC-S4	-2.98	1.79	1.85
2	B	601	59H	CA1-N1	-2.46	1.43	1.46
2	B	601	59H	CX6-CX1	-2.18	1.35	1.39
2	A	601	59H	CA1-N1	-2.08	1.44	1.46
2	A	601	59H	CX6-CX1	-2.08	1.35	1.39
2	B	601	59H	CXA-NX2	2.59	1.43	1.34
2	A	601	59H	CXA-NX2	2.60	1.43	1.34
2	A	601	59H	CX3-CX4	4.24	1.48	1.38
2	B	601	59H	CX3-CX4	4.39	1.48	1.38
2	B	601	59H	CX7-N	4.41	1.45	1.34
2	A	601	59H	C-N1	4.49	1.44	1.34
2	A	601	59H	CX7-N	4.49	1.45	1.34
2	B	601	59H	C-N1	4.64	1.44	1.34
2	A	601	59H	CX5-CX6	6.23	1.51	1.38
2	B	601	59H	CX5-CX6	6.32	1.51	1.38
2	A	601	59H	CX2-CX1	8.66	1.53	1.39
2	B	601	59H	CX2-CX1	8.71	1.53	1.39

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	59H	CX8-NX1-CXA	-4.29	108.23	111.54
2	A	601	59H	CX8-NX1-CXA	-3.99	108.47	111.54
2	B	601	59H	CA3-CXC-S4	-3.87	97.09	103.87
2	A	601	59H	CA3-CXC-S4	-3.30	98.09	103.87
2	A	601	59H	CX8-CX9-NX2	2.08	105.28	102.62
2	A	601	59H	CXC-S4-CXD	2.27	98.79	93.72
2	B	601	59H	CX9-CX8-NX1	2.30	104.82	102.90
2	B	601	59H	CXC-S4-CXD	2.53	99.38	93.72
2	B	601	59H	N-CX7-NX1	2.69	118.22	114.06
2	A	601	59H	CX9-CX8-NX1	3.10	105.48	102.90
2	A	601	59H	N-CX7-NX1	3.10	118.85	114.06
2	B	601	59H	CXD-CA1-N1	3.60	118.57	109.78
2	A	601	59H	CX1-CA-C	4.00	116.75	107.83
2	A	601	59H	CXD-CA1-N1	4.07	119.71	109.78
2	B	601	59H	CX1-CA-C	4.99	118.96	107.83

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	602	GOL	2	0
3	B	602	GOL	2	0
3	B	603	GOL	3	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, 95th 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 ²)	Q<0.9
1	A	515/564 (91%)	0.00	16 (3%) 52 53	15, 33, 69, 111	0
1	B	490/564 (86%)	0.31	39 (7%) 15 16	16, 37, 79, 104	0
All	All	1005/1128 (89%)	0.15	55 (5%) 29 30	15, 34, 76, 111	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	213	ALA	6.6
1	A	50	ALA	6.2
1	A	564	THR	5.6
1	B	192	LYS	4.9
1	B	146	TYR	4.7
1	B	89	LEU	4.5
1	B	211	LYS	4.3
1	B	83	TRP	4.3
1	B	119	GLU	4.2
1	B	122	PHE	4.1
1	B	58	ILE	4.1
1	B	328	TYR	4.1
1	B	139	ALA	4.0
1	B	135	GLU	3.9
1	B	120	ARG	3.8
1	A	199	ARG	3.8
1	B	57	ALA	3.7
1	B	92	ALA	3.6
1	B	212	ASN	3.6
1	B	441	ARG	3.5
1	B	190	PRO	3.5
1	B	143	PRO	3.4
1	B	124	TYR	3.4
1	A	441	ARG	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	563	ALA	3.1
1	A	51	ARG	3.1
1	B	193	ARG	3.1
1	B	84	ALA	3.0
1	A	200	ARG	3.0
1	B	118	ALA	2.9
1	B	560	LEU	2.8
1	B	87	LYS	2.8
1	A	444	ASP	2.8
1	A	560	LEU	2.8
1	A	440	ASP	2.6
1	A	561	PRO	2.6
1	B	189	VAL	2.6
1	B	442	VAL	2.6
1	A	139	ALA	2.6
1	B	91	THR	2.5
1	B	140	LEU	2.5
1	B	59	PRO	2.5
1	B	116	GLN	2.4
1	A	140	LEU	2.3
1	B	327	ARG	2.3
1	A	442	VAL	2.3
1	A	124	TYR	2.3
1	A	395	ARG	2.2
1	B	123	ILE	2.2
1	B	514	THR	2.2
1	B	440	ASP	2.1
1	A	148	ILE	2.0
1	B	90	MET	2.0
1	B	88	GLU	2.0
1	B	102	ALA	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, 95th 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 ²)	Q<0.9
3	GOL	B	602	6/6	0.92	0.19	2.00	29,39,48,48	0
3	GOL	A	602	6/6	0.94	0.16	1.51	21,26,32,38	0
5	IMD	B	607	5/5	0.91	0.12	1.14	48,55,62,63	0
3	GOL	B	603	6/6	0.86	0.17	1.13	47,53,57,59	0
2	59H	B	601	32/32	0.96	0.12	-0.37	28,34,37,41	0
2	59H	A	601	32/32	0.95	0.11	-0.52	23,29,35,38	0
5	IMD	A	604	5/5	0.94	0.07	-	49,52,57,59	0
4	CL	B	605	1/1	0.98	0.07	-	39,39,39,39	0
4	CL	A	603	1/1	0.95	0.09	-	55,55,55,55	0
4	CL	B	606	1/1	0.99	0.08	-	35,35,35,35	0
4	CL	B	604	1/1	0.96	0.06	-	47,47,47,47	0

6.5 Other polymers [i](#)

There are no such residues in this entry.