



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:38 PM GMT

PDB ID : 3UE0
Title : Crystal structure of Acinetobacter baumannii PBP1a in complex with Aztreonam
Authors : Han, S.
Deposited on : 2011-10-28
Resolution : 2.60 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 (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

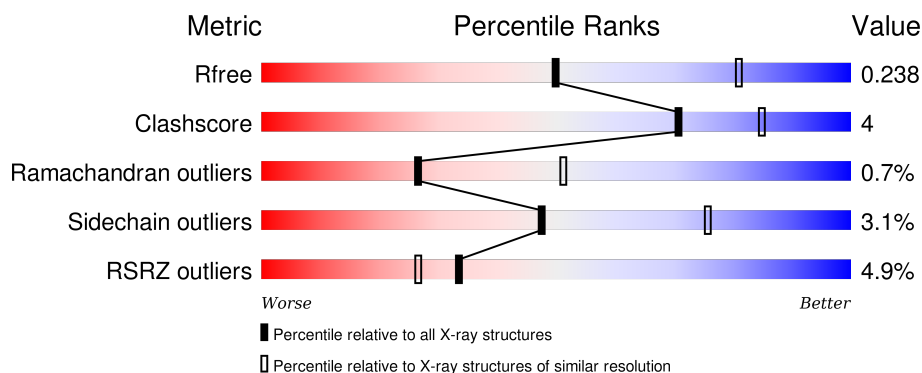
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	2328 (2.60-2.60)
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (2.60-2.60)

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	731	<div> <div>5%</div> <div>71%</div> <div>10%</div> <div>18%</div> </div>
1	B	731	<div> <div>3%</div> <div>73%</div> <div>8%</div> <div>19%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9762 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 Penicillin-binding protein 1a.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	598	Total	C	N	O	S	0	0	0
			4715	3006	832	861	16			
1	B	595	Total	C	N	O	S	0	0	0
			4684	2986	826	856	16			

There are 32 discrepancies between the modelled and reference sequences:

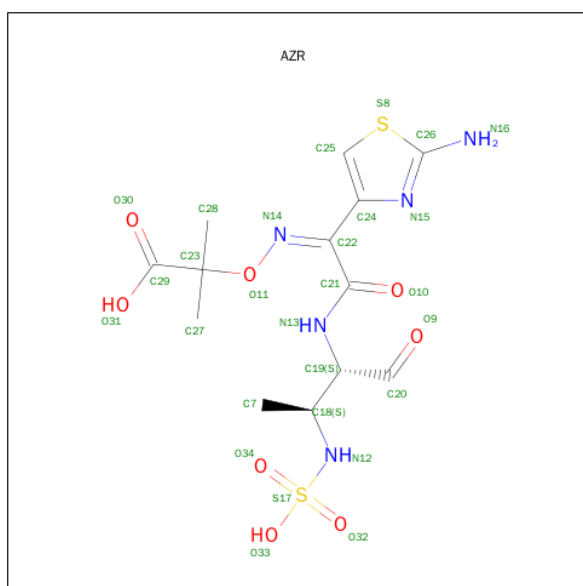
Chain	Residue	Modelled	Actual	Comment	Reference
A	9	MET	-	EXPRESSION TAG	UNP G1C794
A	10	HIS	-	EXPRESSION TAG	UNP G1C794
A	11	HIS	-	EXPRESSION TAG	UNP G1C794
A	12	HIS	-	EXPRESSION TAG	UNP G1C794
A	13	HIS	-	EXPRESSION TAG	UNP G1C794
A	14	HIS	-	EXPRESSION TAG	UNP G1C794
A	15	HIS	-	EXPRESSION TAG	UNP G1C794
A	16	GLU	-	EXPRESSION TAG	UNP G1C794
A	17	ASN	-	EXPRESSION TAG	UNP G1C794
A	18	LEU	-	EXPRESSION TAG	UNP G1C794
A	19	TYR	-	EXPRESSION TAG	UNP G1C794
A	20	PHE	-	EXPRESSION TAG	UNP G1C794
A	21	GLN	-	EXPRESSION TAG	UNP G1C794
A	22	SER	-	EXPRESSION TAG	UNP G1C794
A	23	HIS	-	EXPRESSION TAG	UNP G1C794
A	24	MET	-	EXPRESSION TAG	UNP G1C794
B	9	MET	-	EXPRESSION TAG	UNP G1C794
B	10	HIS	-	EXPRESSION TAG	UNP G1C794
B	11	HIS	-	EXPRESSION TAG	UNP G1C794
B	12	HIS	-	EXPRESSION TAG	UNP G1C794
B	13	HIS	-	EXPRESSION TAG	UNP G1C794
B	14	HIS	-	EXPRESSION TAG	UNP G1C794
B	15	HIS	-	EXPRESSION TAG	UNP G1C794
B	16	GLU	-	EXPRESSION TAG	UNP G1C794
B	17	ASN	-	EXPRESSION TAG	UNP G1C794

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Chain	Residue	Modelled	Actual	Comment	Reference
B	18	LEU	-	EXPRESSION TAG	UNP G1C794
B	19	TYR	-	EXPRESSION TAG	UNP G1C794
B	20	PHE	-	EXPRESSION TAG	UNP G1C794
B	21	GLN	-	EXPRESSION TAG	UNP G1C794
B	22	SER	-	EXPRESSION TAG	UNP G1C794
B	23	HIS	-	EXPRESSION TAG	UNP G1C794
B	24	MET	-	EXPRESSION TAG	UNP G1C794

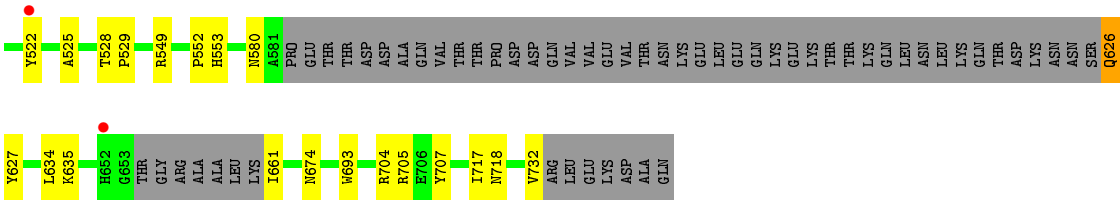
- Molecule 2 is 2-({[(1Z)-1-(2-AMINO-1,3-THIAZOL-4-YL)-2-OXO-2-{[(2S,3S)-1-OXO-3-(SULFOAMINO)BUTAN-2-YL]AMINO}ETHYLIDENE]AMINO}OXY)-2-METHYLPROPANOIC ACID (three-letter code: AZR) (formula: C₁₃H₁₉N₅O₈S₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			28	13	5	8	2		
2	B	1	Total	C	N	O	S	0	0
			28	13	5	8	2		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	127	Total	O	0	0
			127	127		
3	B	180	Total	O	0	0
			180	180		



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2	Depositor
Cell constants a, b, c, α , β , γ	119.08 Å 242.57 Å 49.31 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.94 – 2.60 47.94 – 2.60	Depositor EDS
% Data completeness (in resolution range)	92.6 (47.94-2.60) 92.5 (47.94-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.16 (at 2.61 Å)	Xtriage
Refinement program	BUSTER 2.9.6	Depositor
R, R_{free}	0.191 , 0.241 0.192 , 0.238	Depositor DCC
R_{free} test set	2036 reflections (5.13%)	DCC
Wilson B-factor (Å ²)	42.9	Xtriage
Anisotropy	0.232	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 66.5	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	1 of 41823 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9762	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.13% 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: AZR

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	0.50	0/4824	0.70	1/6541 (0.0%)
1	B	0.52	0/4793	0.71	0/6501
All	All	0.51	0/9617	0.70	1/13042 (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	A	709	GLY	C-N-CA	5.94	136.56	121.70

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	4715	0	4702	48	0
1	B	4684	0	4662	38	0
2	A	28	0	17	6	0
2	B	28	0	16	2	0
3	A	127	0	0	0	0
3	B	180	0	0	3	0
All	All	9762	0	9397	81	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 (81) 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:434:SER:CB	2:A:998:AZR:H20	1.76	1.13
1:A:298:ARG:H	1:A:390:GLN:HE22	1.12	0.95
1:B:298:ARG:H	1:B:390:GLN:HE22	1.10	0.95
1:A:434:SER:OG	2:A:998:AZR:H20	0.76	0.93
1:A:434:SER:HG	2:A:998:AZR:H20	1.14	0.92
1:B:431:GLN:HE21	1:B:528:THR:HA	1.37	0.89
1:A:431:GLN:HE21	1:A:528:THR:HA	1.40	0.85
1:A:549:ARG:NH1	1:A:550:VAL:O	2.15	0.79
1:A:135:LEU:H	1:A:135:LEU:HD12	1.48	0.78
1:A:313:TYR:CE1	1:A:419:GLN:HB2	2.19	0.77
1:B:356:ASN:ND2	1:B:704:ARG:H	1.88	0.71
1:B:434:SER:H	2:B:999:AZR:C20	2.03	0.70
1:A:434:SER:CB	2:A:998:AZR:C20	2.54	0.68
1:B:432:PRO:HB2	1:B:435:THR:HG22	1.79	0.65
1:A:71:PHE:O	1:B:553:HIS:HA	1.99	0.62
1:B:510:ASP:O	1:B:549:ARG:HD3	2.00	0.61
1:A:707:TYR:H	1:A:710:ILE:HG12	1.65	0.61
1:A:528:THR:N	1:A:529:PRO:HD2	2.16	0.60
1:B:528:THR:N	1:B:529:PRO:HD2	2.17	0.60
1:B:465:TRP:HZ3	1:B:522:TYR:HB2	1.67	0.58
1:B:446:GLU:OE2	1:B:635:LYS:HB2	2.04	0.58
1:B:406:GLY:O	1:B:552:PRO:HA	2.04	0.56
1:A:708:GLY:O	1:A:712:ALA:HB3	2.05	0.56
1:A:406:GLY:O	1:A:552:PRO:HA	2.05	0.55
1:A:396:GLY:O	1:A:414:GLY:HA2	2.08	0.53
1:A:137:LYS:NZ	1:B:256:HIS:HD2	2.08	0.52
1:A:380:ASN:ND2	1:A:385:ALA:H	2.07	0.51
1:B:418:TYR:HB2	3:B:873:HOH:O	2.10	0.51
1:A:707:TYR:O	1:A:709:GLY:HA3	2.10	0.50
1:B:37:GLN:HB2	3:B:889:HOH:O	2.10	0.50
1:B:191:ASN:HB3	1:B:194:VAL:HG12	1.93	0.49
1:A:66:ALA:HA	1:A:206:ILE:HG12	1.94	0.49
1:A:553:HIS:HA	1:B:71:PHE:O	2.12	0.49
1:A:323:LYS:HE2	1:A:325:ASN:HD21	1.77	0.49
1:B:356:ASN:HD22	1:B:704:ARG:H	1.59	0.49
1:B:707:TYR:CD2	1:B:707:TYR:N	2.82	0.48
1:B:434:SER:N	2:B:999:AZR:C20	2.76	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:51:GLU:HB2	1:A:54:GLN:HG3	1.96	0.47
1:A:496:LEU:HD22	1:A:525:ALA:HB2	1.96	0.46
1:A:31:VAL:HB	1:A:40:ALA:HB3	1.95	0.46
1:A:69:SER:HB3	3:B:857:HOH:O	2.16	0.46
1:B:323:LYS:HE3	1:B:325:ASN:HD21	1.81	0.45
1:A:432:PRO:HD2	1:A:527:GLY:O	2.15	0.45
1:B:496:LEU:HD22	1:B:525:ALA:HB2	1.97	0.45
1:A:397:GLN:HG3	1:A:693:TRP:HB3	1.99	0.45
1:A:431:GLN:NE2	1:A:528:THR:HA	2.21	0.45
1:B:474:TYR:HE1	1:B:490:THR:HG21	1.81	0.45
1:A:166:ASN:HB2	1:A:234:PRO:HD3	2.00	0.44
1:B:350:ARG:HH11	1:B:358:VAL:HG21	1.82	0.44
1:A:514:GLN:HB3	1:A:517:GLN:HG2	1.99	0.44
1:A:474:TYR:HE1	1:A:490:THR:HG21	1.83	0.44
1:B:397:GLN:HG3	1:B:693:TRP:HB3	1.98	0.44
1:A:640:TYR:HD1	1:A:732:VAL:HG23	1.82	0.44
1:A:135:LEU:H	1:A:135:LEU:CD1	2.24	0.44
1:A:247:GLU:HG2	1:A:417:PHE:HE1	1.81	0.44
1:A:255:LYS:HG3	1:B:185:LYS:HE3	2.00	0.43
1:A:29:LEU:HD22	1:A:266:TYR:O	2.18	0.43
1:B:626:GLN:OE1	1:B:626:GLN:N	2.52	0.43
1:B:66:ALA:HA	1:B:206:ILE:HG12	2.01	0.43
1:A:72:PHE:O	1:B:553:HIS:HB2	2.19	0.42
1:B:381:GLU:CD	1:B:381:GLU:H	2.22	0.42
1:B:661:ILE:HD11	1:B:717:ILE:HA	2.01	0.42
1:B:528:THR:H	1:B:529:PRO:HD2	1.85	0.42
1:A:350:ARG:HH11	1:A:358:VAL:HG21	1.84	0.42
1:A:511:PHE:HA	1:A:549:ARG:HD3	2.02	0.42
1:B:221:TYR:O	1:B:225:VAL:HG22	2.20	0.42
1:B:707:TYR:N	1:B:707:TYR:HD2	2.17	0.42
1:A:221:TYR:O	1:A:225:VAL:HG22	2.20	0.41
1:B:431:GLN:NE2	1:B:528:THR:HA	2.19	0.41
1:A:519:PRO:HD2	1:A:524:ILE:HG22	2.02	0.41
1:A:434:SER:HB2	2:A:998:AZR:C20	2.44	0.41
1:A:298:ARG:H	1:A:390:GLN:NE2	1.96	0.41
1:B:432:PRO:HD3	1:B:529:PRO:O	2.20	0.41
1:A:29:LEU:HB3	1:A:42:TYR:HB2	2.03	0.41
1:A:674:ASN:N	2:A:998:AZR:N15	2.69	0.41
1:A:626:GLN:HA	1:A:627:TYR:HA	1.84	0.41
1:B:166:ASN:HB2	1:B:234:PRO:HD3	2.02	0.41
1:B:443:LEU:O	1:B:446:GLU:HB2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:29:LEU:HB3	1:B:42:TYR:HB2	2.03	0.41
1:A:380:ASN:HD21	1:A:385:ALA:H	1.68	0.40
1:A:247:GLU:HG2	1:A:417:PHE:CE1	2.55	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	590/731 (81%)	557 (94%)	27 (5%)	6 (1%)	19	39
1	B	587/731 (80%)	562 (96%)	23 (4%)	2 (0%)	46	72
All	All	1177/1462 (80%)	1119 (95%)	50 (4%)	8 (1%)	26	51

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	383	LYS
1	A	708	GLY
1	B	627	TYR
1	A	349	ALA
1	B	349	ALA
1	A	384	THR
1	A	463	GLY
1	A	709	GLY

5.3.2 Protein sidechains [i](#)

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	490/608 (81%)	473 (96%)	17 (4%)	43	71
1	B	486/608 (80%)	473 (97%)	13 (3%)	52	79
All	All	976/1216 (80%)	946 (97%)	30 (3%)	47	76

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	35	ASP
1	A	59	PHE
1	A	135	LEU
1	A	139	ASP
1	A	145	VAL
1	A	198	ARG
1	A	218	GLN
1	A	353	ARG
1	A	381	GLU
1	A	384	THR
1	A	421	LYS
1	A	514	GLN
1	A	674	ASN
1	A	698	GLN
1	A	710	ILE
1	A	718	ASN
1	A	732	VAL
1	B	35	ASP
1	B	59	PHE
1	B	145	VAL
1	B	218	GLN
1	B	435	THR
1	B	514	GLN
1	B	580	ASN
1	B	626	GLN
1	B	634	LEU
1	B	674	ASN
1	B	705	ARG
1	B	718	ASN
1	B	732	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	61	HIS
1	A	222	GLN
1	A	256	HIS
1	A	325	ASN
1	A	380	ASN
1	A	390	GLN
1	A	431	GLN
1	A	514	GLN
1	A	556	GLN
1	A	629	GLN
1	A	652	HIS
1	B	256	HIS
1	B	325	ASN
1	B	356	ASN
1	B	390	GLN
1	B	431	GLN
1	B	514	GLN
1	B	556	GLN
1	B	629	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](#)

2 ligands 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	$\# Z > 2$
2	AZR	A	998	1	18,28,28	1.32	3 (16%)	17,41,41	2.05	3 (17%)
2	AZR	B	999	1	18,28,28	1.92	6 (33%)	17,41,41	2.44	7 (41%)

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	AZR	A	998	1	-	0/20/35/35	0/1/1/1
2	AZR	B	999	1	-	0/20/35/35	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	999	AZR	O11-N14	-4.24	1.34	1.42
2	B	999	AZR	C24-C22	-3.02	1.43	1.48
2	B	999	AZR	C19-C18	-2.70	1.51	1.54
2	B	999	AZR	C19-N13	-2.46	1.43	1.46
2	A	998	AZR	C19-C18	-2.13	1.52	1.54
2	A	998	AZR	O34-S17	2.26	1.44	1.42
2	B	999	AZR	O34-S17	2.90	1.45	1.42
2	A	998	AZR	O32-S17	3.02	1.45	1.42
2	B	999	AZR	O32-S17	3.67	1.45	1.42

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	999	AZR	O32-S17-N12	-5.77	100.79	108.50
2	A	998	AZR	C27-C23-C29	-5.62	98.63	111.64
2	B	999	AZR	C27-C23-C29	-4.41	101.44	111.64
2	B	999	AZR	C18-C19-C20	-2.89	106.31	112.19
2	B	999	AZR	O10-C21-C22	-2.48	117.76	120.53
2	B	999	AZR	C28-C23-C29	-2.26	106.40	111.64
2	B	999	AZR	N16-C26-N15	2.91	126.73	122.92
2	A	998	AZR	N16-C26-N15	3.04	126.90	122.92
2	B	999	AZR	O10-C21-N13	3.30	128.94	123.01
2	A	998	AZR	O10-C21-N13	3.71	129.68	123.01

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	998	AZR	6	0
2	B	999	AZR	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, 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	598/731 (81%)	0.18	39 (6%)	22 16	23, 54, 97, 123	0
1	B	595/731 (81%)	-0.01	20 (3%)	49 41	15, 44, 93, 143	0
All	All	1193/1462 (81%)	0.08	59 (4%)	33 26	15, 49, 95, 143	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	460	ILE	7.9
1	B	465	TRP	7.2
1	A	363	SER	6.4
1	A	362	PRO	5.3
1	A	327	SER	5.2
1	B	463	GLY	5.2
1	B	464	LYS	5.0
1	B	466	THR	4.9
1	A	562	TYR	4.8
1	B	462	ILE	4.8
1	B	459	PRO	4.5
1	A	328	SER	4.3
1	A	382	ALA	4.3
1	A	371	VAL	4.1
1	A	364	ARG	3.9
1	A	326	SER	3.9
1	A	324	VAL	3.6
1	B	226	ALA	3.5
1	A	329	PHE	3.4
1	B	469	ASN	3.4
1	B	522	TYR	3.4
1	A	565	VAL	3.3
1	A	368	ILE	3.3
1	A	347	SER	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	342	GLN	3.2
1	B	470	SER	3.2
1	B	225	VAL	3.1
1	A	704	ARG	2.9
1	A	345	GLY	2.9
1	A	463	GLY	2.9
1	B	228	PRO	2.8
1	A	343	TRP	2.8
1	B	471	ASP	2.8
1	A	626	GLN	2.7
1	B	198	ARG	2.7
1	B	190	TYR	2.7
1	B	467	PRO	2.6
1	A	346	MET	2.6
1	B	652	HIS	2.5
1	A	341	VAL	2.4
1	A	357	SER	2.4
1	A	321	VAL	2.4
1	A	356	ASN	2.4
1	A	323	LYS	2.4
1	A	325	ASN	2.3
1	A	355	ALA	2.3
1	A	361	ALA	2.3
1	B	461	THR	2.2
1	A	464	LYS	2.2
1	A	190	TYR	2.2
1	A	734	LEU	2.2
1	A	661	ILE	2.1
1	A	372	LYS	2.1
1	A	373	ASP	2.1
1	A	344	SER	2.1
1	A	369	VAL	2.1
1	B	194	VAL	2.1
1	A	381	GLU	2.0
1	A	733	ARG	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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(\AA^2)	Q<0.9
2	AZR	A	998	28/28	0.91	0.24	1.72	61,80,93,93	0
2	AZR	B	999	28/28	0.90	0.21	0.61	54,65,92,96	0

6.5 Other polymers [i](#)

There are no such residues in this entry.