Full wwPDB X-ray Structure Validation Report (i)

Dec 6, 2016 – 03:18 PM EST

PDB ID	:	5HLB
Title	:	E. coli PBP1b in complex with acyl-aztreonam and moenomycin
Authors	:	King, D.T.; Strynadka, N.C.J.
Deposited on	:	2016-01-14
Resolution	:	2.42 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

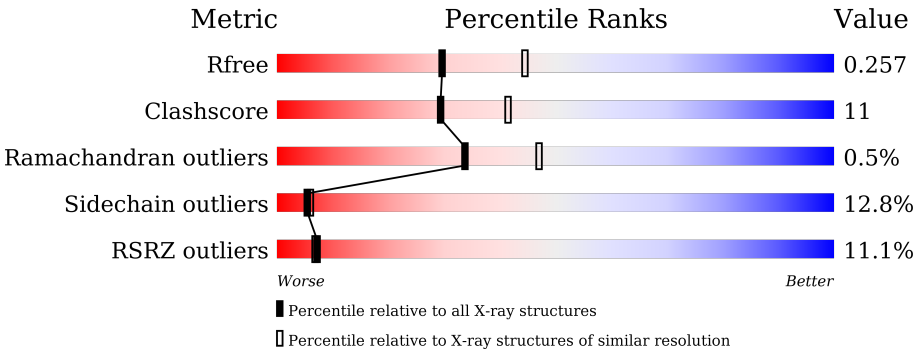
MolProbity	:	4.02b-467
Mogul	:	1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20028442
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20028442

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.42 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	3386 (2.44-2.40)
Clashscore	102246	3897 (2.44-2.40)
Ramachandran outliers	100387	3837 (2.44-2.40)
Sidechain outliers	100360	3838 (2.44-2.40)
RSRZ outliers	91569	3396 (2.44-2.40)

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 >=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 <=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	747	11% 74% 20%

2 Entry composition

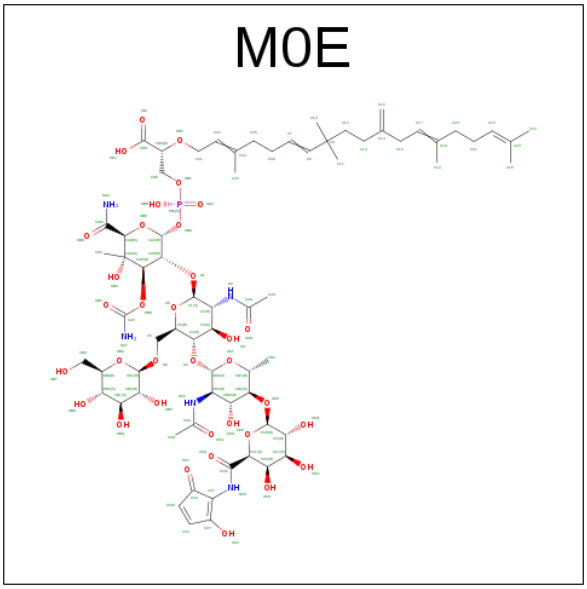
There are 4 unique types of molecules in this entry. The entry contains 6076 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 1B.

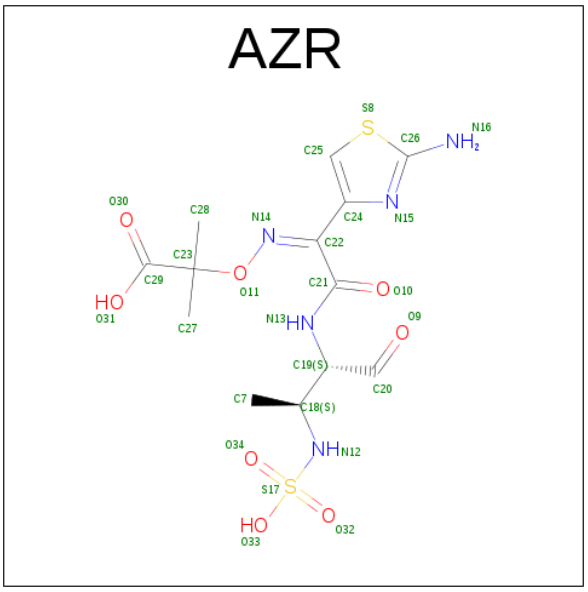
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	733	Total	C	N	O	S	0	0	0
			5738	3638	1012	1062	26			

- Molecule 2 is MOENOMYCIN (three-letter code: M0E) (formula: $C_{69}H_{106}N_5O_{34}P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			77	39	5	32	1		

- Molecule 3 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: $C_{13}H_{19}N_5O_8S_2$).



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

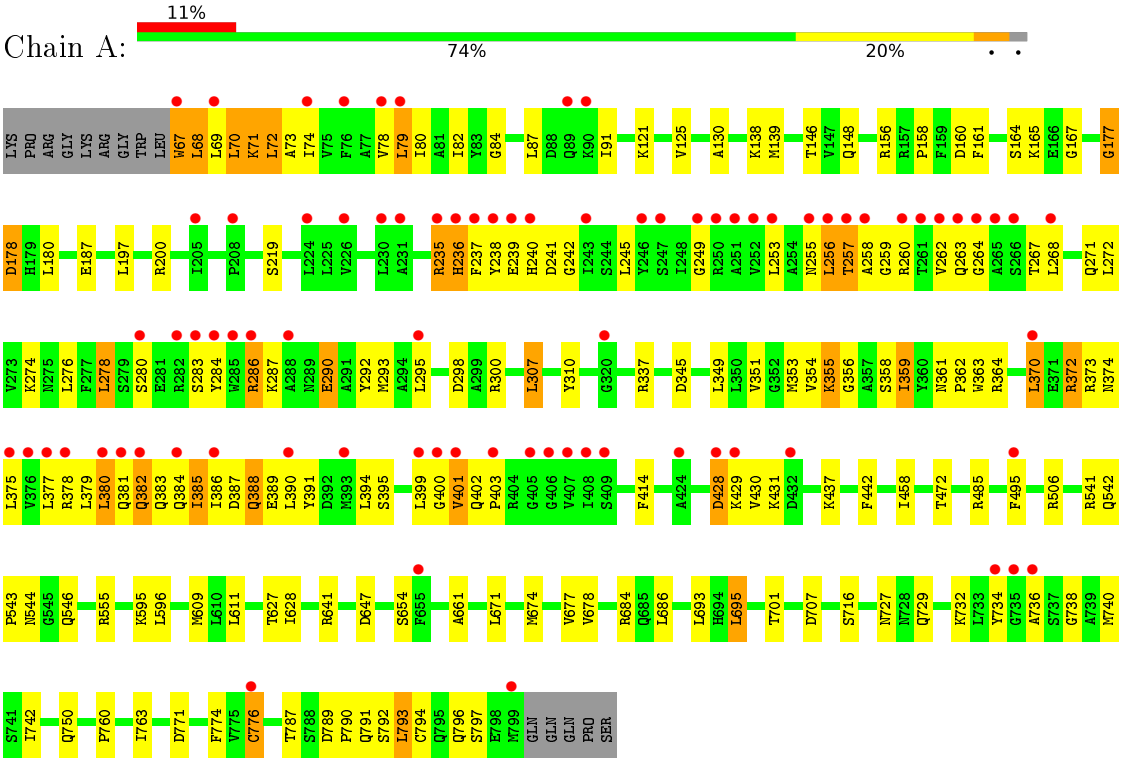
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	233	Total	O	0	0
			233	233		

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 (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: Penicillin-binding protein 1B



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, α, β, γ	62.53Å 64.39Å 301.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.43 – 2.42 62.97 – 2.42	Depositor EDS
% Data completeness (in resolution range)	96.6 (35.43-2.42) 89.4 (62.97-2.42)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.24 (at 2.42Å)	Xtriage
Refinement program	BUSTER 2.10.0	Depositor
R, R_{free}	0.225 , 0.243 0.238 , 0.257	Depositor DCC
R_{free} test set	2377 reflections (5.46%)	DCC
Wilson B-factor (Å^2)	36.0	Xtriage
Anisotropy	0.760	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.29 , 34.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.41$, $< L^2 > = 0.23$	Xtriage
Estimated twinning fraction	0.076 for k,h,-l	Xtriage
F_o,F_c correlation	0.90	EDS
Total number of atoms	6076	wwPDB-VP
Average B, all atoms (Å^2)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.48% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 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, M0E

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	0.47	0/5854	0.77	2/7947 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	384	GLN	N-CA-C	7.16	130.34	111.00
1	A	257	THR	N-CA-C	-6.75	92.77	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	237	PHE	Peptide

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	5738	0	5792	119	0
2	A	77	0	57	20	0
3	A	28	0	16	1	0
4	A	233	0	0	0	0
All	All	6076	0	5865	132	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 11.

All (132) 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)
2:A:901:M0E:CAX	2:A:901:M0E:OBG	1.64	1.45
2:A:901:M0E:O1	2:A:901:M0E:CAR	1.65	1.41
1:A:372:ARG:NH1	1:A:375:LEU:HD22	1.36	1.35
1:A:258:ALA:HB1	1:A:259:GLY:HA3	1.21	1.18
1:A:355:LYS:HD2	1:A:372:ARG:NH2	1.60	1.15
1:A:355:LYS:HD2	1:A:372:ARG:HH21	0.96	1.05
1:A:372:ARG:HH12	1:A:375:LEU:CD2	1.76	0.99
2:A:901:M0E:CAR	2:A:901:M0E:C1	2.41	0.98
1:A:71:LYS:HG3	1:A:72:LEU:N	1.78	0.95
1:A:71:LYS:HG3	1:A:72:LEU:H	1.32	0.94
1:A:68:LEU:HG	1:A:71:LYS:HG2	1.49	0.94
1:A:69:LEU:O	1:A:73:ALA:CB	2.17	0.93
1:A:400:GLY:HA3	1:A:402:GLN:HG2	1.50	0.92
1:A:69:LEU:O	1:A:73:ALA:HB2	1.69	0.92
1:A:355:LYS:CD	1:A:372:ARG:HH21	1.84	0.91
1:A:374:ASN:HA	1:A:377:LEU:HD12	1.54	0.88
1:A:258:ALA:CB	1:A:259:GLY:HA3	1.97	0.86
1:A:372:ARG:HH11	1:A:375:LEU:HD22	1.40	0.85
1:A:372:ARG:NH1	1:A:375:LEU:CD2	2.30	0.84
1:A:355:LYS:O	2:A:901:M0E:H5	1.78	0.83
1:A:258:ALA:HB1	1:A:259:GLY:CA	2.10	0.80
1:A:372:ARG:HH12	1:A:375:LEU:HD22	0.98	0.80
1:A:388:GLN:HA	1:A:391:TYR:HB3	1.64	0.80
1:A:362:PRO:HB3	1:A:399:LEU:HD13	1.65	0.79
1:A:262:VAL:CG1	1:A:264:GLY:H	1.95	0.79
2:A:901:M0E:O1	2:A:901:M0E:CAP	2.32	0.78
1:A:262:VAL:HG11	2:A:901:M0E:HAU2	1.50	0.76
1:A:240:HIS:CD2	1:A:240:HIS:H	2.07	0.72
1:A:79:LEU:HA	1:A:82:ILE:HG12	1.72	0.70
1:A:283:SER:HB3	1:A:286:ARG:HB2	1.73	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:283:SER:CB	1:A:286:ARG:HB2	2.22	0.69
1:A:71:LYS:O	1:A:74:ILE:HG22	1.92	0.69
1:A:69:LEU:O	1:A:73:ALA:HB3	1.91	0.69
1:A:363:TRP:CZ3	1:A:402:GLN:HB3	2.29	0.68
2:A:901:M0E:PBI	2:A:901:M0E:CAX	2.82	0.67
2:A:901:M0E:O3	2:A:901:M0E:OCF	2.13	0.64
1:A:91:ILE:HD11	1:A:293:MET:HG2	1.80	0.64
1:A:262:VAL:HG13	1:A:264:GLY:H	1.61	0.63
1:A:67:TRP:CE3	1:A:67:TRP:HA	2.32	0.63
1:A:506:ARG:HD2	1:A:707:ASP:OD1	2.00	0.61
2:A:901:M0E:HAS2	2:A:901:M0E:OBD	1.99	0.61
1:A:239:GLU:CB	1:A:298:ASP:HB3	2.30	0.61
2:A:901:M0E:OBG	2:A:901:M0E:CAR	2.45	0.61
1:A:388:GLN:HA	1:A:391:TYR:CB	2.31	0.61
1:A:377:LEU:HD22	1:A:394:LEU:HB3	1.82	0.60
1:A:399:LEU:HD12	1:A:400:GLY:N	2.18	0.58
1:A:359:ILE:HD12	2:A:901:M0E:HCB1	1.86	0.57
1:A:776:CYS:HA	1:A:793:LEU:HD13	1.86	0.57
2:A:901:M0E:HO3	2:A:901:M0E:CBU	2.13	0.56
1:A:239:GLU:HB2	1:A:298:ASP:HB3	1.87	0.56
1:A:262:VAL:HG13	1:A:264:GLY:N	2.21	0.55
1:A:380:LEU:HD23	1:A:386:ILE:HG12	1.87	0.55
1:A:379:LEU:HD23	1:A:382:GLN:HE22	1.73	0.54
1:A:349:LEU:O	1:A:353:MET:HG2	2.08	0.54
1:A:67:TRP:HE3	1:A:67:TRP:HA	1.73	0.53
1:A:458:ILE:HD11	1:A:472:THR:HB	1.89	0.53
1:A:235:ARG:HG2	1:A:238:TYR:OH	2.08	0.53
1:A:262:VAL:HG13	1:A:263:GLN:N	2.24	0.53
1:A:377:LEU:CD2	1:A:394:LEU:HB3	2.38	0.52
1:A:760:PRO:HD2	1:A:763:ILE:HG13	1.90	0.52
1:A:239:GLU:N	1:A:240:HIS:O	2.43	0.52
1:A:283:SER:HB2	1:A:286:ARG:HB2	1.91	0.52
1:A:736:ALA:O	1:A:740:MET:HB2	2.10	0.52
1:A:370:LEU:HG	1:A:399:LEU:HD23	1.90	0.52
1:A:701:THR:HG21	1:A:734:TYR:HA	1.92	0.51
1:A:283:SER:HB3	1:A:286:ARG:H	1.75	0.51
1:A:139:MET:HG3	1:A:146:THR:HG23	1.93	0.51
1:A:789:ASP:C	1:A:791:GLN:H	2.15	0.51
1:A:349:LEU:HB2	1:A:373:ARG:HG3	1.93	0.50
1:A:240:HIS:H	1:A:240:HIS:HD2	1.59	0.50
1:A:70:LEU:O	1:A:74:ILE:N	2.42	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:596:LEU:HA	1:A:627:THR:HG21	1.94	0.50
1:A:278:LEU:HD11	1:A:290:GLU:HB2	1.93	0.50
1:A:595:LYS:HB3	1:A:661:ALA:HB1	1.93	0.50
1:A:354:VAL:HG22	2:A:901:M0E:HAH3	1.93	0.50
1:A:611:LEU:O	3:A:902:AZR:H25	2.11	0.49
1:A:239:GLU:N	1:A:240:HIS:C	2.65	0.49
1:A:355:LYS:O	2:A:901:M0E:H1	2.13	0.49
1:A:161:PHE:HB2	1:A:164:SER:O	2.12	0.48
1:A:241:ASP:HB3	1:A:267:THR:HG21	1.95	0.48
2:A:901:M0E:HAH2	2:A:901:M0E:OBC	2.13	0.48
1:A:262:VAL:CG1	1:A:264:GLY:N	2.73	0.48
1:A:789:ASP:O	1:A:791:GLN:N	2.46	0.48
1:A:240:HIS:CD2	1:A:240:HIS:N	2.80	0.48
1:A:262:VAL:CG1	2:A:901:M0E:HAU2	2.25	0.48
1:A:274:LYS:HA	1:A:278:LEU:HD12	1.94	0.48
1:A:236:HIS:HA	1:A:238:TYR:CD2	2.48	0.47
1:A:363:TRP:HH2	1:A:403:PRO:HD2	1.79	0.47
1:A:370:LEU:CD2	1:A:373:ARG:HH21	2.27	0.47
1:A:674:MET:HA	1:A:677:VAL:HG13	1.96	0.47
1:A:160:ASP:HA	1:A:165:LYS:HG3	1.97	0.47
1:A:249:GLY:O	1:A:262:VAL:O	2.33	0.47
2:A:901:M0E:HAQ	2:A:901:M0E:OBG	2.15	0.47
1:A:130:ALA:O	1:A:200:ARG:HD3	2.14	0.46
1:A:428:ASP:O	1:A:430:VAL:N	2.40	0.46
1:A:307:LEU:HA	1:A:310:TYR:HB3	1.98	0.46
1:A:239:GLU:CA	1:A:240:HIS:C	2.84	0.46
2:A:901:M0E:HAH2	2:A:901:M0E:CAV	2.46	0.45
1:A:271:GLN:HG2	2:A:901:M0E:HAT2	1.80	0.45
1:A:353:MET:O	1:A:356:GLY:O	2.35	0.45
1:A:693:LEU:HB3	1:A:695:LEU:HD13	1.98	0.45
1:A:378:ARG:O	1:A:381:GLN:HG2	2.17	0.45
1:A:84:GLY:HA2	1:A:87:LEU:HD12	1.98	0.45
1:A:377:LEU:C	1:A:391:TYR:HE2	2.20	0.44
1:A:776:CYS:CA	1:A:793:LEU:HD13	2.47	0.44
1:A:274:LYS:HA	1:A:278:LEU:CD1	2.48	0.44
1:A:241:ASP:CB	1:A:267:THR:HG21	2.48	0.43
1:A:242:GLY:O	1:A:245:LEU:N	2.50	0.43
1:A:363:TRP:CH2	1:A:403:PRO:HD2	2.53	0.43
1:A:375:LEU:O	1:A:378:ARG:HG3	2.18	0.43
1:A:727:ASN:OD1	1:A:729:GLN:HB2	2.19	0.43
1:A:177:GLY:HA2	1:A:178:ASP:HA	1.87	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:268:LEU:HD23	1:A:271:GLN:OE1	2.18	0.43
1:A:414:PHE:CE1	1:A:442:PHE:HB2	2.54	0.43
1:A:283:SER:HB3	1:A:286:ARG:N	2.34	0.42
1:A:385:ILE:HD13	1:A:385:ILE:H	1.84	0.42
1:A:774:PHE:N	1:A:796:GLN:HG2	2.34	0.42
1:A:345:ASP:HB3	1:A:394:LEU:HD22	2.02	0.42
1:A:377:LEU:HB3	1:A:391:TYR:CE2	2.55	0.42
1:A:262:VAL:HG12	1:A:264:GLY:H	1.80	0.42
1:A:738:GLY:O	1:A:742:ILE:HG12	2.20	0.42
1:A:431:LYS:HZ3	1:A:431:LYS:HB3	1.85	0.41
1:A:256:LEU:O	1:A:258:ALA:N	2.53	0.41
1:A:378:ARG:HB3	1:A:391:TYR:OH	2.20	0.41
1:A:362:PRO:HB3	1:A:399:LEU:CD1	2.42	0.41
1:A:678:VAL:O	1:A:684:ARG:HA	2.21	0.41
2:A:901:M0E:CAQ	2:A:901:M0E:OBG	2.69	0.41
1:A:351:VAL:O	1:A:354:VAL:HG12	2.20	0.41
1:A:158:PRO:HB3	1:A:167:GLY:HA2	2.03	0.41
1:A:278:LEU:HD13	1:A:286:ARG:HG2	2.03	0.40
1:A:121:LYS:O	1:A:125:VAL:HG23	2.20	0.40
1:A:542:GLN:HB2	1:A:543:PRO:CD	2.52	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	731/747 (98%)	677 (93%)	50 (7%)	4 (0%)	34 47

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	790	PRO

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Mol	Chain	Res	Type
1	A	177	GLY
1	A	78	VAL
1	A	401	VAL

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	615/627 (98%)	536 (87%)	79 (13%)	5 6

All (79) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	67	TRP
1	A	68	LEU
1	A	70	LEU
1	A	71	LYS
1	A	72	LEU
1	A	79	LEU
1	A	80	ILE
1	A	138	LYS
1	A	148	GLN
1	A	156	ARG
1	A	178	ASP
1	A	180	LEU
1	A	187	GLU
1	A	197	LEU
1	A	219	SER
1	A	235	ARG
1	A	236	HIS
1	A	253	LEU
1	A	255	ASN
1	A	256	LEU
1	A	257	THR
1	A	260	ARG
1	A	272	LEU

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Mol	Chain	Res	Type
1	A	276	LEU
1	A	278	LEU
1	A	280	SER
1	A	284	TYR
1	A	286	ARG
1	A	287	LYS
1	A	290	GLU
1	A	292	TYR
1	A	295	LEU
1	A	300	ARG
1	A	307	LEU
1	A	337	ARG
1	A	355	LYS
1	A	358	SER
1	A	359	ILE
1	A	361	ASN
1	A	364	ARG
1	A	370	LEU
1	A	372	ARG
1	A	380	LEU
1	A	382	GLN
1	A	383	GLN
1	A	385	ILE
1	A	387	ASP
1	A	388	GLN
1	A	389	GLU
1	A	390	LEU
1	A	395	SER
1	A	401	VAL
1	A	428	ASP
1	A	429	LYS
1	A	437	LYS
1	A	485	ARG
1	A	495	PHE
1	A	541	ARG
1	A	544	ASN
1	A	546	GLN
1	A	555	ARG
1	A	609	MET
1	A	628	ILE
1	A	641	ARG
1	A	647	ASP

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Mol	Chain	Res	Type
1	A	654	SER
1	A	671	LEU
1	A	686	LEU
1	A	695	LEU
1	A	716	SER
1	A	732	LYS
1	A	750	GLN
1	A	771	ASP
1	A	776	CYS
1	A	787	THR
1	A	792	SER
1	A	793	LEU
1	A	794	CYS
1	A	797	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	189	ASN
1	A	240	HIS
1	A	382	GLN
1	A	542	GLN

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

There are no carbohydrates in this entry.

5.6 Ligand geometry

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	M0E	A	901	-	75,81,114	3.15	34 (45%)	95,122,166	3.42	49 (51%)
3	AZR	A	902	1	19,28,28	3.12	7 (36%)	16,41,41	2.61	5 (31%)

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	M0E	A	901	-	-	2/47/158/206	0/5/5/6
3	AZR	A	902	1	-	0/21/35/35	0/1/1/1

All (41) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	901	M0E	ODF-CDG	-10.16	1.18	1.42
3	A	902	AZR	C25-S8	-9.72	1.55	1.70
2	A	901	M0E	CAR-CAP	-6.53	1.38	1.52
2	A	901	M0E	OBR-CBM	-5.89	1.29	1.43
2	A	901	M0E	OCQ-CCM	-4.49	1.15	1.23
2	A	901	M0E	CAO-CAP	-4.11	1.47	1.53
3	A	902	AZR	C19-N13	-3.91	1.42	1.46
2	A	901	M0E	CBK-CBL	-3.67	1.42	1.52
2	A	901	M0E	C6-C5	-3.16	1.41	1.51
2	A	901	M0E	CCM-NCS	-3.09	1.27	1.32
2	A	901	M0E	CBW-CBV	-3.09	1.46	1.53
2	A	901	M0E	PBI-OBB	-3.01	1.42	1.55
2	A	901	M0E	OCG-CCA	-2.96	1.16	1.23
3	A	902	AZR	O11-N14	-2.93	1.36	1.43
2	A	901	M0E	CBO-CBN	-2.65	1.42	1.51
2	A	901	M0E	OBQ-CBL	-2.55	1.36	1.43
2	A	901	M0E	OBC-CAV	-2.55	1.17	1.21
2	A	901	M0E	CBW-CBX	-2.43	1.45	1.52
2	A	901	M0E	CAQ-CAW	-2.15	1.45	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	901	M0E	CCJ-CCI	-2.11	1.46	1.52
2	A	901	M0E	OCR-CCK	-2.01	1.38	1.43
2	A	901	M0E	CBJ-CBK	2.06	1.58	1.52
2	A	901	M0E	OCE-CBX	2.09	1.49	1.43
2	A	901	M0E	PBI-OBG	2.10	1.66	1.60
2	A	901	M0E	PBI-OBF	2.26	1.68	1.59
3	A	902	AZR	S17-N12	2.29	1.62	1.60
3	A	902	AZR	C26-N16	2.51	1.42	1.35
2	A	901	M0E	CBZ-CBY	2.61	1.57	1.51
2	A	901	M0E	O5-C5	2.63	1.50	1.44
2	A	901	M0E	CBV-NCC	2.75	1.50	1.45
2	A	901	M0E	CCA-NCC	2.89	1.45	1.34
2	A	901	M0E	O6-CBJ	3.05	1.45	1.40
3	A	902	AZR	C21-N13	3.12	1.40	1.34
2	A	901	M0E	OCP-CCL	3.74	1.49	1.43
2	A	901	M0E	O6-C6	3.80	1.50	1.43
2	A	901	M0E	OBG-CAX	4.44	1.64	1.42
2	A	901	M0E	CAH-CAG	4.54	1.60	1.50
3	A	902	AZR	C24-N15	5.55	1.54	1.37
2	A	901	M0E	OCD-CBW	7.03	1.59	1.43
2	A	901	M0E	O1-CAR	8.45	1.65	1.43
2	A	901	M0E	CAW-NAU	11.64	1.55	1.32

All (54) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	901	M0E	C1-C2-N2	-10.17	91.87	111.01
2	A	901	M0E	CBL-CBM-CBN	-8.85	94.45	110.23
2	A	901	M0E	O3-C3-C4	-8.67	89.35	109.89
2	A	901	M0E	CBJ-CBK-CBL	-7.79	94.53	109.98
2	A	901	M0E	CBZ-CBY-CBX	-7.39	101.22	113.38
2	A	901	M0E	OBE-CAX-OBG	-6.93	102.22	111.36
3	A	902	AZR	O34-S17-O32	-6.79	101.75	120.14
2	A	901	M0E	OBT-CBO-CBN	-6.05	91.09	111.30
2	A	901	M0E	C1-O5-C5	-5.91	102.14	113.74
2	A	901	M0E	C1-O1-CAR	-5.57	103.20	118.00
2	A	901	M0E	O5-C5-C4	-5.04	99.02	109.78
2	A	901	M0E	ODF-CDG-CDK	-4.65	95.92	110.16
2	A	901	M0E	OCO-CCJ-CCK	-4.63	99.93	110.36
2	A	901	M0E	OCQ-CCM-NCS	-4.56	115.50	123.06
2	A	901	M0E	OBS-CBN-CBO	-4.44	94.89	106.38
2	A	901	M0E	OCD-CBW-CBX	-4.11	100.14	109.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	902	AZR	C24-C25-S8	-4.03	106.84	111.79
2	A	901	M0E	CAS-CAO-CAP	-4.00	107.72	111.69
2	A	901	M0E	C3-C4-C5	-3.78	102.20	110.85
2	A	901	M0E	OBD-CAW-NAU	-3.58	117.14	123.06
2	A	901	M0E	OBP-CBK-CBL	-3.46	102.56	110.36
2	A	901	M0E	C6-O6-CBJ	-2.91	107.73	113.81
2	A	901	M0E	OCR-CCK-CCJ	-2.79	104.06	110.36
2	A	901	M0E	OBG-CAX-CAR	-2.79	103.19	108.39
3	A	902	AZR	N16-C26-N15	-2.72	119.36	122.92
2	A	901	M0E	O1-CAR-CAP	-2.65	99.68	106.83
2	A	901	M0E	ODF-CDG-CDH	-2.50	104.08	111.70
2	A	901	M0E	CBO-CBN-CBM	-2.36	107.06	112.99
2	A	901	M0E	CBU-CBV-NCC	-2.34	106.60	111.01
2	A	901	M0E	OCG-CCA-NCC	-2.21	117.33	121.84
3	A	902	AZR	O9-C20-C19	-2.18	119.07	125.81
2	A	901	M0E	OCN-CCI-CCH	-2.12	105.30	110.01
2	A	901	M0E	OBC-CAV-NAT	-2.08	121.68	125.50
2	A	901	M0E	OBB-PBI-OAZ	-2.04	101.93	112.56
2	A	901	M0E	CCH-OCE-CBX	-2.03	112.62	118.00
2	A	901	M0E	OBA-CAO-CAP	2.04	112.94	107.72
2	A	901	M0E	C6-C5-C4	2.09	118.38	113.33
2	A	901	M0E	CBK-CBL-CBM	2.12	114.70	110.79
2	A	901	M0E	OCF-CBY-CBZ	2.20	111.25	106.57
2	A	901	M0E	O6-CBJ-CBK	2.23	110.74	108.00
2	A	901	M0E	O5-C5-C6	2.62	112.02	106.61
2	A	901	M0E	OCF-CBY-CBX	2.76	114.15	109.09
2	A	901	M0E	O4-C4-C5	2.81	116.82	109.33
2	A	901	M0E	OCR-CCK-CCL	3.34	117.12	109.80
2	A	901	M0E	OBR-CBM-CBL	3.36	117.93	110.36
2	A	901	M0E	CCB-CCA-NCC	3.48	122.77	116.10
2	A	901	M0E	OBG-PBI-OAZ	3.69	124.45	109.34
2	A	901	M0E	OBQ-CBL-CBM	3.84	119.02	110.36
2	A	901	M0E	OBS-CBJ-CBK	3.87	118.33	110.28
2	A	901	M0E	CAQ-CAW-NAU	4.16	121.81	117.12
3	A	902	AZR	O32-S17-N12	5.08	115.69	108.63
2	A	901	M0E	C3-C2-N2	5.61	122.32	110.67
2	A	901	M0E	CCL-CCM-NCS	7.06	125.76	116.97
2	A	901	M0E	OBS-CBN-CBM	7.63	124.23	109.67

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	901	M0E	CAP-OBH-CAV-OBC
2	A	901	M0E	CAP-OBH-CAV-NAT

There are no ring outliers.

2 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	901	M0E	20	0
3	A	902	AZR	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^{th} 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	733/747 (98%)	0.47	81 (11%) 7 7	16, 46, 122, 172	0

All (81) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	246	TYR	12.5
1	A	401	VAL	10.0
1	A	283	SER	8.1
1	A	408	ILE	8.0
1	A	256	LEU	7.5
1	A	282	ARG	7.2
1	A	260	ARG	7.1
1	A	247	SER	6.9
1	A	257	THR	6.8
1	A	237	PHE	6.5
1	A	240	HIS	6.4
1	A	734	TYR	6.2
1	A	258	ALA	5.9
1	A	262	VAL	5.8
1	A	284	TYR	5.7
1	A	67	TRP	5.6
1	A	261	THR	5.6
1	A	735	GLY	5.5
1	A	285	TRP	5.5
1	A	280	SER	5.3
1	A	406	GLY	5.2
1	A	390	LEU	5.1
1	A	286	ARG	4.7
1	A	255	ASN	4.6
1	A	252	VAL	4.6
1	A	238	TYR	4.4
1	A	89	GLN	4.1

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Mol	Chain	Res	Type	RSRZ
1	A	243	ILE	4.1
1	A	382	GLN	4.0
1	A	253	LEU	4.0
1	A	264	GLY	3.8
1	A	378	ARG	3.7
1	A	405	GLY	3.7
1	A	265	ALA	3.7
1	A	78	VAL	3.6
1	A	386	ILE	3.5
1	A	495	PHE	3.5
1	A	263	GLN	3.4
1	A	69	LEU	3.4
1	A	799	MET	3.3
1	A	432	ASP	3.3
1	A	409	SER	3.3
1	A	428	ASP	3.2
1	A	249	GLY	3.2
1	A	407	VAL	3.2
1	A	230	LEU	3.2
1	A	205	ILE	3.2
1	A	90	LYS	3.2
1	A	736	ALA	3.0
1	A	429	LYS	2.9
1	A	380	LEU	2.8
1	A	251	ALA	2.8
1	A	320	GLY	2.8
1	A	384	GLN	2.7
1	A	76	PHE	2.7
1	A	224	LEU	2.6
1	A	403	PRO	2.6
1	A	295	LEU	2.6
1	A	266	SER	2.5
1	A	375	LEU	2.5
1	A	79	LEU	2.5
1	A	208	PRO	2.4
1	A	399	LEU	2.4
1	A	370	LEU	2.4
1	A	393	MET	2.4
1	A	239	GLU	2.3
1	A	288	ALA	2.3
1	A	377	LEU	2.3
1	A	376	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	381	GLN	2.3
1	A	268	LEU	2.3
1	A	776	CYS	2.2
1	A	400	GLY	2.2
1	A	226	VAL	2.1
1	A	655	PHE	2.1
1	A	231	ALA	2.1
1	A	250	ARG	2.1
1	A	236	HIS	2.0
1	A	74	ILE	2.0
1	A	235	ARG	2.0
1	A	424	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, 95^{th} 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	AZR	A	902	28/28	0.96	0.16	-0.20	33,47,73,76	0
2	M0E	A	901	77/109	0.85	0.16	-0.78	79,106,125,132	0

6.5 Other polymers (i)

There are no such residues in this entry.