



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 11:12 PM GMT

PDB ID : 5I0P
Title : Crystal Structure of a Beta-lactamase domain protein from Burkholderia ambifaria
Authors : Seattle Structural Genomics Center for Infectious Disease (SSGCID)
Deposited on : 2016-02-04
Resolution : 2.50 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 : unknown
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

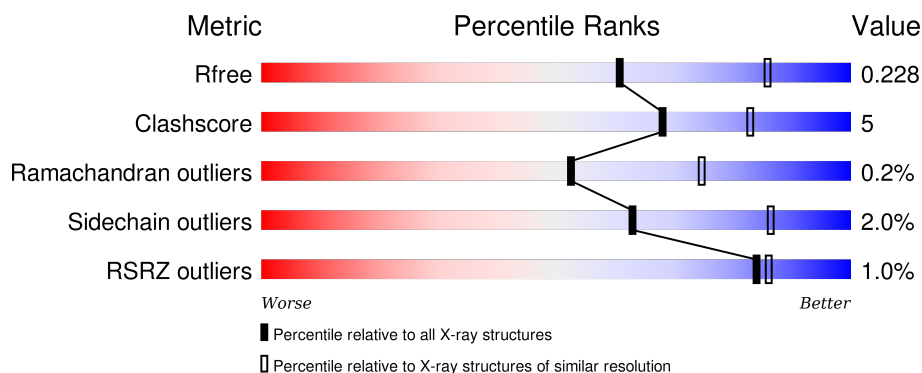
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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	366	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 12%, green 75%, grey 12%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 75% 12% • 12% </div> </div>
1	B	366	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 9%, green 79%, grey 12%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 79% 9% • 12% </div> </div>
1	C	366	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 9%, green 83%, grey 7%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 83% 9% • 7% </div> </div>
1	D	366	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 12%, green 81%, grey 6%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 81% 12% • 6% </div> </div>

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
2	ZN	D	401	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10562 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 Beta-lactamase domain protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	322	Total	C	N	O	S	0	1	0
			2485	1577	451	443	14			
1	B	323	Total	C	N	O	S	0	2	0
			2466	1571	447	434	14			
1	C	339	Total	C	N	O	S	0	1	0
			2602	1657	472	457	16			
1	D	344	Total	C	N	O	S	0	0	0
			2677	1699	488	475	15			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	MET	-	expression tag	UNP B1YNR3
A	-6	ALA	-	expression tag	UNP B1YNR3
A	-5	HIS	-	expression tag	UNP B1YNR3
A	-4	HIS	-	expression tag	UNP B1YNR3
A	-3	HIS	-	expression tag	UNP B1YNR3
A	-2	HIS	-	expression tag	UNP B1YNR3
A	-1	HIS	-	expression tag	UNP B1YNR3
A	0	HIS	-	expression tag	UNP B1YNR3
B	-7	MET	-	expression tag	UNP B1YNR3
B	-6	ALA	-	expression tag	UNP B1YNR3
B	-5	HIS	-	expression tag	UNP B1YNR3
B	-4	HIS	-	expression tag	UNP B1YNR3
B	-3	HIS	-	expression tag	UNP B1YNR3
B	-2	HIS	-	expression tag	UNP B1YNR3
B	-1	HIS	-	expression tag	UNP B1YNR3
B	0	HIS	-	expression tag	UNP B1YNR3
C	-7	MET	-	expression tag	UNP B1YNR3
C	-6	ALA	-	expression tag	UNP B1YNR3
C	-5	HIS	-	expression tag	UNP B1YNR3
C	-4	HIS	-	expression tag	UNP B1YNR3
C	-3	HIS	-	expression tag	UNP B1YNR3

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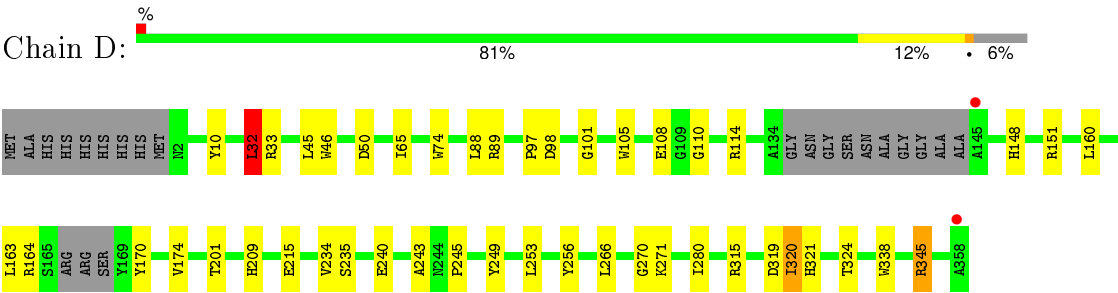
Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	HIS	-	expression tag	UNP B1YNR3
C	-1	HIS	-	expression tag	UNP B1YNR3
C	0	HIS	-	expression tag	UNP B1YNR3
D	-7	MET	-	expression tag	UNP B1YNR3
D	-6	ALA	-	expression tag	UNP B1YNR3
D	-5	HIS	-	expression tag	UNP B1YNR3
D	-4	HIS	-	expression tag	UNP B1YNR3
D	-3	HIS	-	expression tag	UNP B1YNR3
D	-2	HIS	-	expression tag	UNP B1YNR3
D	-1	HIS	-	expression tag	UNP B1YNR3
D	0	HIS	-	expression tag	UNP B1YNR3

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	2	Total Zn 2 2	0	0
2	A	2	Total Zn 2 2	0	0
2	D	2	Total Zn 2 2	0	0
2	C	2	Total Zn 2 2	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	59	Total O 59 59	0	0
3	B	49	Total O 49 49	0	0
3	C	100	Total O 100 100	0	0
3	D	116	Total O 116 116	0	0



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	45.17Å 141.23Å 236.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.94 – 2.50 44.93 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.8 (44.94-2.50) 99.9 (44.93-2.50)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.03 (at 2.51Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.166 , 0.229 0.164 , 0.228	Depositor DCC
R_{free} test set	1996 reflections (3.72%)	DCC
Wilson B-factor (Å ²)	36.0	Xtriage
Anisotropy	0.647	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 43.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 53685 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	10562	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.95% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.55	1/2554 (0.0%)	0.74	5/3480 (0.1%)
1	B	0.54	0/2539	0.71	1/3466 (0.0%)
1	C	0.58	0/2676	0.77	4/3648 (0.1%)
1	D	0.57	0/2748	0.73	3/3743 (0.1%)
All	All	0.56	1/10517 (0.0%)	0.74	13/14337 (0.1%)

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	C	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	107	CYS	CB-SG	-5.20	1.73	1.81

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	32	LEU	CA-CB-CG	9.00	136.01	115.30
1	A	32	LEU	CA-CB-CG	8.12	133.97	115.30
1	C	33	ARG	NE-CZ-NH2	-6.47	117.07	120.30
1	A	224	ASP	CB-CG-OD1	6.06	123.76	118.30
1	D	32	LEU	CB-CG-CD2	-6.02	100.76	111.00
1	A	183	ARG	NE-CZ-NH1	-5.99	117.30	120.30
1	A	32	LEU	CB-CG-CD2	-5.76	101.21	111.00
1	C	284	ARG	NE-CZ-NH2	-5.69	117.45	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	33	ARG	NE-CZ-NH2	-5.63	117.48	120.30
1	B	224	ASP	CB-CG-OD1	5.59	123.33	118.30
1	C	32	LEU	CA-CB-CG	5.59	128.16	115.30
1	D	89	ARG	NE-CZ-NH2	-5.42	117.59	120.30
1	C	284	ARG	CG-CD-NE	-5.04	101.22	111.80

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	146	ALA	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	2485	0	2382	27	0
1	B	2466	0	2339	20	0
1	C	2602	0	2481	21	0
1	D	2677	0	2586	27	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
3	A	59	0	0	3	0
3	B	49	0	0	1	0
3	C	100	0	0	2	0
3	D	116	0	0	2	0
All	All	10562	0	9788	95	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 5.

All (95) 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:257:GLU:HG3	1:B:280:ILE:HG21	1.69	0.75
1:C:338:TRP:CE2	1:C:345:ARG:HD2	2.22	0.74
1:D:32:LEU:HD22	1:D:45:LEU:HB2	1.71	0.71
1:C:250:LEU:O	1:C:284:ARG:NH2	2.24	0.71
1:D:319:ASP:O	1:D:321:HIS:N	2.24	0.71
1:B:234:VAL:HG13	1:B:245:PRO:HB2	1.72	0.70
1:A:338:TRP:CE2	1:A:345:ARG:HD2	2.28	0.69
1:A:108:GLU:OE1	3:A:501:HOH:O	2.11	0.68
1:C:256:TYR:HB2	1:C:280:ILE:HD11	1.76	0.67
1:C:108:GLU:OE2	3:C:501:HOH:O	2.13	0.66
1:B:40:LEU:HD21	1:B:269:HIS:ND1	2.11	0.66
1:D:256:TYR:HB2	1:D:280:ILE:CD1	2.29	0.62
1:C:185:ARG:NE	3:C:503:HOH:O	2.23	0.61
1:C:338:TRP:CZ2	1:C:345:ARG:HD2	2.35	0.61
1:C:33:ARG:NH2	1:C:270:GLY:O	2.35	0.59
1:C:66:SER:OG	1:C:71:ARG:NH2	2.32	0.59
1:A:40:LEU:HD21	1:A:269:HIS:CE1	2.38	0.58
1:C:234:VAL:HG13	1:C:245:PRO:HB2	1.86	0.57
1:D:148:HIS:HA	1:D:151:ARG:NH2	2.20	0.57
1:A:30:ARG:NH1	1:A:81:HIS:O	2.37	0.57
1:A:234:VAL:HG13	1:A:245:PRO:HB2	1.88	0.56
1:A:189:ALA:HB2	1:A:198:ARG:NH2	2.21	0.56
1:A:32:LEU:HD11	1:A:47:LEU:HG	1.89	0.55
1:D:234:VAL:HG13	1:D:245:PRO:HB2	1.88	0.54
1:D:249:TYR:CZ	1:D:253:LEU:HD21	2.43	0.54
1:A:338:TRP:CZ2	1:A:345:ARG:HD2	2.42	0.54
1:A:32:LEU:HD22	1:A:45:LEU:HB2	1.90	0.54
1:D:256:TYR:HB2	1:D:280:ILE:HD11	1.88	0.54
1:D:32:LEU:HD21	1:D:74:TRP:CH2	2.44	0.53
1:D:215:GLU:HG3	3:D:544:HOH:O	2.08	0.53
1:B:338:TRP:CZ2	1:B:345:ARG:HD3	2.44	0.53
1:D:65:ILE:HG23	1:D:98:ASP:HA	1.90	0.52
1:D:97:PRO:HG3	1:D:170:TYR:CD1	2.44	0.52
1:A:187:GLY:O	1:A:198:ARG:NH1	2.43	0.51
1:C:40:LEU:HD21	1:C:269:HIS:CE1	2.45	0.50
1:B:338:TRP:CE2	1:B:345:ARG:HD2	2.48	0.49
1:C:119:LEU:HD22	1:C:178:PRO:HG2	1.94	0.49
1:A:236:VAL:HB	1:A:331:LEU:HB3	1.95	0.49
1:B:119:LEU:HD22	1:B:178:PRO:HG2	1.94	0.49
1:B:201:THR:HA	1:B:209:HIS:O	2.13	0.49
1:B:297:ALA:HB1	1:B:311:ILE:HD13	1.94	0.49
1:C:87:VAL:HB	1:C:115:TRP:CE2	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:160:LEU:HB3	1:D:164:ARG:HH21	1.79	0.48
1:D:201:THR:HA	1:D:209:HIS:O	2.14	0.48
1:C:65:ILE:HG23	1:C:98:ASP:HA	1.96	0.47
1:D:338:TRP:CD2	1:D:345:ARG:HD3	2.50	0.47
1:A:131:LEU:HD12	1:A:131:LEU:HA	1.55	0.47
1:A:101:GLY:HA2	1:A:174:VAL:HG22	1.98	0.46
1:A:201:THR:HA	1:A:209:HIS:O	2.15	0.46
1:D:105:TRP:CE2	1:D:110:GLY:HA2	2.50	0.46
1:A:257:GLU:HG3	1:A:280:ILE:HG21	1.98	0.46
1:D:114:ARG:HD3	3:D:527:HOH:O	2.16	0.45
1:B:33:ARG:NH2	1:B:270:GLY:O	2.49	0.45
1:C:96:HIS:CE1	1:C:205:HIS:NE2	2.85	0.45
1:A:19:ALA:HB1	1:A:73:HIS:CD2	2.51	0.45
1:D:10:TYR:CD1	1:D:271:LYS:HE3	2.52	0.45
1:C:256:TYR:HB2	1:C:280:ILE:CD1	2.45	0.45
1:A:32:LEU:HD21	1:A:74:TRP:CH2	2.52	0.45
1:D:98:ASP:N	1:D:98:ASP:OD2	2.49	0.45
1:A:148:HIS:HA	1:A:151:ARG:NH1	2.32	0.44
1:C:46:TRP:CH2	1:C:266:LEU:HB3	2.51	0.44
1:D:33:ARG:NH2	1:D:270:GLY:O	2.50	0.44
1:A:224:ASP:HA	1:A:227:LEU:HD21	1.99	0.43
1:B:40:LEU:HD21	1:B:269:HIS:CE1	2.54	0.43
1:B:102:LEU:HD23	1:B:102:LEU:HA	1.84	0.43
1:B:46:TRP:HB2	1:B:61:VAL:HB	2.01	0.43
1:D:101:GLY:HA2	1:D:174:VAL:HG22	2.00	0.43
1:B:105:TRP:CE2	1:B:110:GLY:HA2	2.54	0.43
1:B:40:LEU:HD21	1:B:269:HIS:HD1	1.83	0.43
1:D:240:GLU:CD	1:D:243:ALA:HB2	2.39	0.43
1:B:338:TRP:CE2	1:B:345:ARG:CD	3.02	0.43
1:D:50:ASP:HB3	1:D:88:LEU:HD12	2.00	0.43
1:A:123:LEU:O	1:A:127:MET:HB2	2.19	0.42
1:A:105:TRP:CE2	1:A:110:GLY:HA2	2.54	0.42
1:D:320:ILE:O	1:D:324:THR:HG23	2.20	0.42
1:B:21:ASP:OD1	1:B:22:THR:N	2.50	0.42
1:B:308:ILE:HA	1:B:308:ILE:HD13	1.91	0.42
1:D:163:LEU:HA	1:D:163:LEU:HD23	1.85	0.42
1:C:98:ASP:N	1:C:98:ASP:OD2	2.52	0.42
1:D:164:ARG:HG3	1:D:164:ARG:H	1.55	0.41
1:B:120:TRP:CE3	1:B:182[B]:ARG:HD2	2.55	0.41
1:C:201:THR:HA	1:C:209:HIS:O	2.20	0.41
1:D:235:SER:O	1:D:245:PRO:HG3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:40:LEU:HD21	1:C:269:HIS:ND1	2.35	0.41
1:A:103:ALA:HB3	3:A:530:HOH:O	2.20	0.41
1:D:46:TRP:CH2	1:D:266:LEU:HB3	2.56	0.41
1:A:252:SER:HB3	3:A:536:HOH:O	2.20	0.41
1:B:218:GLY:HA2	3:B:541:HOH:O	2.21	0.41
1:A:308:ILE:HD13	1:A:308:ILE:HA	1.86	0.41
1:B:19:ALA:HB1	1:B:73:HIS:CD2	2.56	0.40
1:C:259:MET:O	1:C:277:ARG:NH2	2.51	0.40
1:C:350:ASP:CG	1:C:354:ARG:HH22	2.24	0.40
1:A:51:GLU:HA	1:A:55:GLN:O	2.22	0.40
1:A:102:LEU:HD23	1:A:102:LEU:HA	1.80	0.40
1:A:331:LEU:HA	1:A:331:LEU:HD23	1.83	0.40

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	315/366 (86%)	310 (98%)	5 (2%)	0	100	100
1	B	317/366 (87%)	312 (98%)	5 (2%)	0	100	100
1	C	334/366 (91%)	325 (97%)	8 (2%)	1 (0%)	46	68
1	D	338/366 (92%)	332 (98%)	5 (2%)	1 (0%)	46	68
All	All	1304/1464 (89%)	1279 (98%)	23 (2%)	2 (0%)	52	75

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	320	ILE
1	C	315	ARG

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	248/295 (84%)	242 (98%)	6 (2%)	57	82
1	B	241/295 (82%)	234 (97%)	7 (3%)	50	77
1	C	255/295 (86%)	251 (98%)	4 (2%)	70	90
1	D	270/295 (92%)	266 (98%)	4 (2%)	72	91
All	All	1014/1180 (86%)	993 (98%)	21 (2%)	63	85

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	LEU
1	A	39	SER
1	A	127	MET
1	A	171	SER
1	A	290	ARG
1	A	331	LEU
1	B	169	TYR
1	B	177	VAL
1	B	182[A]	ARG
1	B	182[B]	ARG
1	B	235	SER
1	B	301	LYS
1	B	345	ARG
1	C	32	LEU
1	C	151	ARG
1	C	171	SER
1	C	284	ARG
1	D	32	LEU
1	D	108	GLU
1	D	315	ARG
1	D	345	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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](#)

Of 8 ligands modelled in this entry, 8 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	322/366 (87%)	-0.40	2 (0%) 90 91	21, 37, 67, 90	0
1	B	323/366 (88%)	-0.17	5 (1%) 76 79	25, 45, 75, 106	0
1	C	339/366 (92%)	-0.28	4 (1%) 81 83	19, 31, 66, 112	0
1	D	344/366 (93%)	-0.36	2 (0%) 90 91	20, 33, 63, 84	0
All	All	1328/1464 (90%)	-0.30	13 (0%) 84 86	19, 36, 69, 112	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	315	ARG	4.0
1	C	144	ALA	3.7
1	A	312	MET	3.1
1	B	351	GLY	2.8
1	D	358	ALA	2.7
1	B	302	PRO	2.7
1	C	311	ILE	2.6
1	D	145	ALA	2.5
1	A	326	ALA	2.5
1	C	314	ARG	2.2
1	B	341	GLY	2.2
1	B	349	ASP	2.0
1	B	150	ALA	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	ZN	D	401	1/1	0.98	0.15	2.01	38,38,38,38	0
2	ZN	C	401	1/1	0.99	0.11	-1.00	36,36,36,36	0
2	ZN	A	401	1/1	0.99	0.10	-1.11	37,37,37,37	0
2	ZN	B	401	1/1	0.98	0.08	-2.24	38,38,38,38	0
2	ZN	C	400	1/1	1.00	0.06	-3.42	43,43,43,43	0
2	ZN	A	400	1/1	0.98	0.06	-3.89	50,50,50,50	0
2	ZN	D	400	1/1	0.98	0.08	-3.98	47,47,47,47	0
2	ZN	B	400	1/1	0.98	0.04	-4.97	57,57,57,57	0

6.5 Other polymers [i](#)

There are no such residues in this entry.