



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 07:15 AM UTC

PDB ID : 4OLD / pdb_00004old
Title : Crystal structure of AmpC beta-lactamase in complex with the product form of (6R,7R)-7-amino-8-oxo-5-thia-1-azabicyclo[4.2.0]oct-2-ene-2-carboxylic acid
Authors : Shoichet, B.K.; Barelier, S.
Deposited on : 2014-01-23
Resolution : 1.48 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

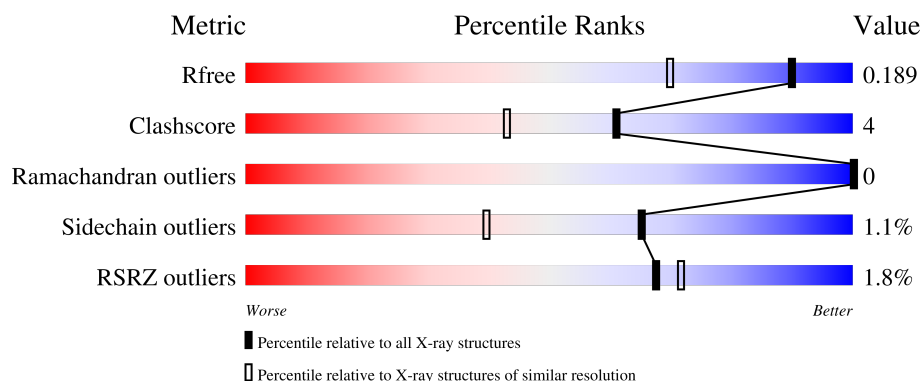
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

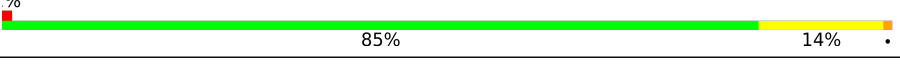
The reported resolution of this entry is 1.48 Å.

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}	180053	6779 (1.50-1.46)
Clashscore	190562	7025 (1.50-1.46)
Ramachandran outliers	187476	6917 (1.50-1.46)
Sidechain outliers	187428	6914 (1.50-1.46)
RSRZ outliers	180081	6781 (1.50-1.46)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	358	 3% 78% 18% ..
1	B	358	 % 85% 14% .

2 Entry composition [i](#)

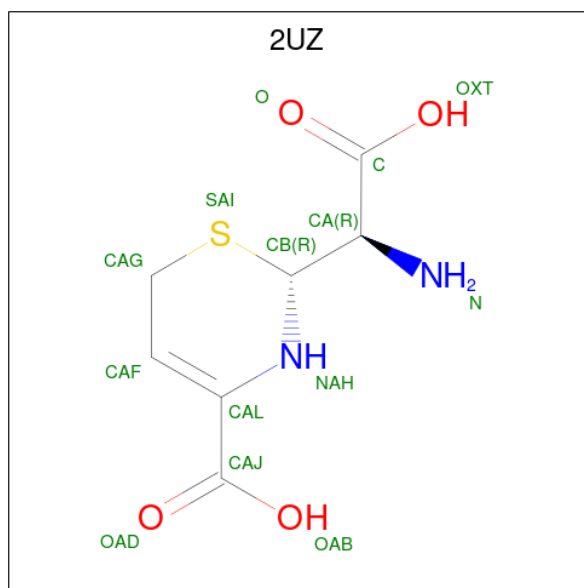
There are 4 unique types of molecules in this entry. The entry contains 6386 atoms, of which 0 are hydrogens and 0 are deuteriums.

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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	351	Total	C	N	O	S	0	16	0
			2834	1823	479	524	8			
1	B	358	Total	C	N	O	S	0	12	0
			2878	1848	489	534	7			

- Molecule 2 is (2R)-2-[(R)-amino(carboxy)methyl]-3,6-dihydro-2H-1,3-thiazine-4-carboxylic acid (CCD ID: 2UZ) (formula: C₇H₁₀N₂O₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	1
			28	14	4	8	2		
2	B	1	Total	C	N	O	S	0	0
			14	7	2	4	1		

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	1
			5	4	1		
3	A	1	Total	O	P	0	0
			5	4	1		
3	B	1	Total	O	P	0	1
			5	4	1		

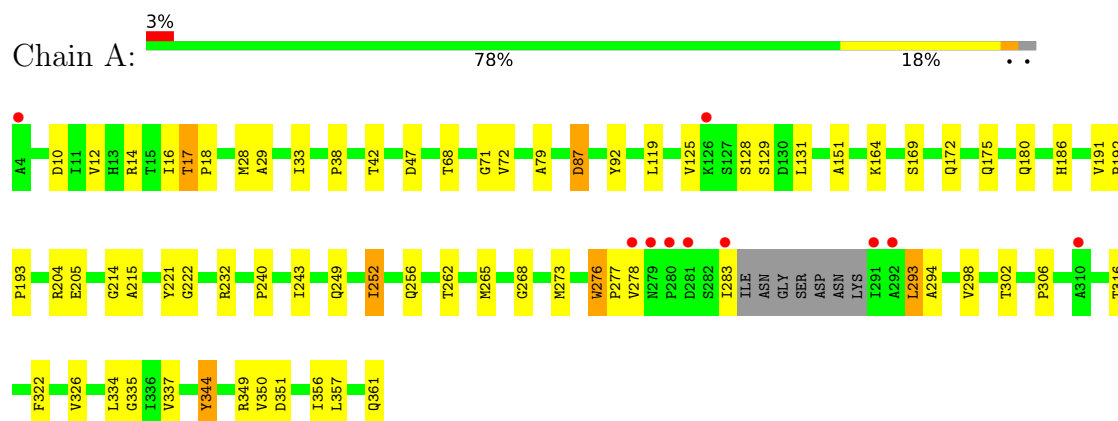
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	292	Total	O	0	15
			292	292		
4	B	325	Total	O	0	13
			325	325		

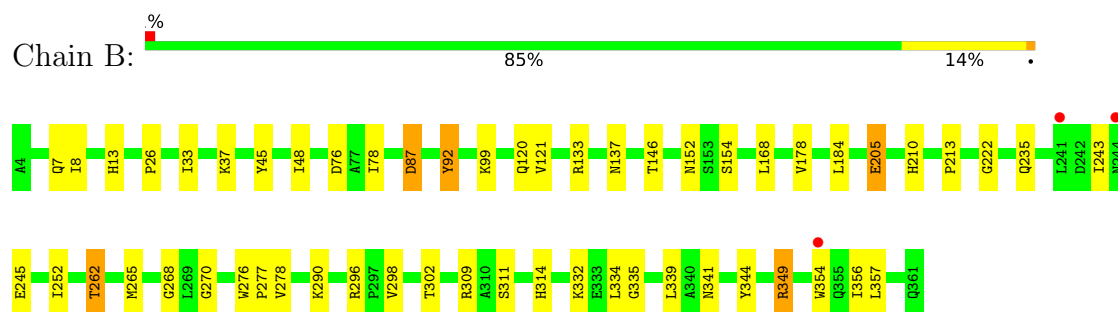
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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: Beta-lactamase



• Molecule 1: Beta-lactamase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	118.45Å 76.42Å 97.88Å 90.00° 115.95° 90.00°	Depositor
Resolution (Å)	29.34 – 1.48 29.34 – 1.48	Depositor EDS
% Data completeness (in resolution range)	99.3 (29.34-1.48) 99.3 (29.34-1.48)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.01 (at 1.48Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.164 , 0.188 (Not available) , 0.189	Depositor DCC
R_{free} test set	4891 reflections (3.78%)	wwPDB-VP
Wilson B-factor (Å ²)	16.3	Xtriage
Anisotropy	0.020	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 41.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	6386	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.26% 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, 0.333 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: 2UZ, PO4

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	1.75	31/2913 (1.1%)	1.49	25/3985 (0.6%)
1	B	1.69	28/2958 (0.9%)	1.39	7/4046 (0.2%)
All	All	1.72	59/5871 (1.0%)	1.44	32/8031 (0.4%)

All (59) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	92	TYR	CZ-OH	7.96	1.54	1.38
1	A	337	VAL	C-O	7.59	1.31	1.24
1	A	186	HIS	ND1-CE1	7.27	1.39	1.32
1	A	42	THR	C-O	6.82	1.32	1.23
1	A	38	PRO	C-O	6.68	1.30	1.23
1	B	45	TYR	C-O	-6.66	1.16	1.24
1	B	76	ASP	C-O	6.65	1.32	1.24
1	A	298	VAL	C-O	-6.62	1.17	1.24
1	A	87	ASP	CA-C	6.11	1.59	1.53
1	B	87	ASP	CG-OD2	-6.09	1.13	1.25
1	B	245	GLU	C-O	6.05	1.31	1.24
1	B	37	LYS	CA-C	6.01	1.61	1.52
1	B	133	ARG	C-O	-6.00	1.17	1.24
1	A	356	ILE	CA-CB	5.94	1.61	1.54
1	B	349	ARG	N-CA	5.93	1.53	1.46
1	A	186	HIS	CG-ND1	-5.87	1.31	1.38
1	A	276	TRP	NE1-CE2	-5.83	1.31	1.37
1	A	79	ALA	C-O	5.77	1.31	1.24
1	A	335	GLY	C-O	5.75	1.29	1.23
1	B	222	GLY	N-CA	5.72	1.52	1.45
1	B	13	HIS	CG-ND1	-5.72	1.31	1.38
1	A	47	ASP	C-O	-5.70	1.17	1.24
1	A	273	MET	N-CA	5.59	1.52	1.45
1	A	222	GLY	N-CA	5.59	1.52	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	335	GLY	N-CA	5.57	1.51	1.45
1	A	164	LYS	N-CA	5.55	1.51	1.46
1	A	344	TYR	C-O	-5.54	1.17	1.23
1	A	326	VAL	C-O	5.51	1.29	1.24
1	B	178	VAL	C-O	-5.46	1.17	1.24
1	A	128	SER	CA-C	5.46	1.59	1.52
1	A	17	THR	C-N	5.44	1.40	1.33
1	A	268	GLY	N-CA	5.43	1.51	1.45
1	A	215	ALA	CA-C	5.41	1.59	1.52
1	A	221	TYR	C-O	5.41	1.27	1.23
1	B	341	ASN	C-O	-5.40	1.15	1.23
1	B	121	VAL	CA-C	5.40	1.57	1.52
1	A	71	GLY	C-O	-5.38	1.17	1.23
1	A	16	ILE	CA-CB	5.36	1.61	1.54
1	B	235	GLN	C-O	5.34	1.30	1.24
1	B	87	ASP	CB-CG	5.33	1.65	1.52
1	B	213	PRO	C-O	5.30	1.29	1.23
1	A	322	PHE	C-O	-5.29	1.17	1.23
1	A	306	PRO	CA-CB	5.26	1.61	1.53
1	B	184	LEU	CA-CB	5.26	1.59	1.52
1	B	356	ILE	C-O	5.20	1.29	1.24
1	B	268	GLY	C-O	5.20	1.28	1.24
1	B	154	SER	C-O	-5.17	1.18	1.24
1	B	339	LEU	N-CA	5.17	1.52	1.46
1	A	42	THR	CA-C	-5.14	1.46	1.52
1	B	137	ASN	N-CA	5.12	1.52	1.46
1	A	294	ALA	CA-C	5.11	1.59	1.53
1	B	314	HIS	CD2-NE2	5.09	1.43	1.37
1	A	186	HIS	N-CA	-5.07	1.39	1.46
1	B	146	THR	N-CA	5.05	1.53	1.46
1	A	193	PRO	CA-CB	5.02	1.61	1.53
1	A	192	PRO	N-CA	5.01	1.53	1.46
1	B	311	SER	CA-C	-5.01	1.46	1.52
1	B	210	HIS	ND1-CE1	5.01	1.37	1.32
1	B	99	LYS	N-CA	-5.00	1.39	1.46

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	192	PRO	O-C-N	-7.60	117.58	121.15
1	B	302	THR	CA-C-O	-6.83	114.10	120.04
1	A	169	SER	N-CA-C	-6.49	101.93	110.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	302	THR	O-C-N	6.33	127.54	120.83
1	A	172	GLN	CA-C-O	6.22	127.15	120.55
1	A	356	ILE	N-CA-C	6.15	116.38	110.47
1	A	12	VAL	O-C-N	-6.07	115.59	121.90
1	A	214	GLY	O-C-N	-6.04	117.94	123.56
1	A	10	ASP	N-CA-C	6.01	117.63	111.14
1	A	302	THR	CA-C-O	-5.87	114.94	120.04
1	A	14	ARG	CA-CB-CG	-5.82	102.47	114.10
1	B	78	ILE	N-CA-C	-5.76	104.89	110.42
1	A	72	VAL	CA-C-O	-5.74	114.76	120.85
1	A	205	GLU	N-CA-C	-5.72	104.02	111.74
1	A	172	GLN	O-C-N	-5.68	116.09	122.12
1	A	214	GLY	CA-C-O	5.66	127.54	121.37
1	A	232	ARG	CG-CD-NE	-5.52	99.85	112.00
1	A	172	GLN	N-CA-C	5.48	117.25	111.28
1	A	252	ILE	O-C-N	-5.44	116.58	121.91
1	A	191	VAL	O-C-N	5.36	127.21	121.10
1	B	168	LEU	CA-C-O	-5.31	115.36	121.47
1	B	121	VAL	N-CA-C	-5.25	104.10	108.99
1	A	129[A]	SER	CA-C-O	-5.24	115.32	120.82
1	A	129[B]	SER	CA-C-O	-5.24	115.32	120.82
1	B	332	LYS	CA-CB-CG	5.18	124.47	114.10
1	A	68	THR	CA-C-O	-5.17	114.03	119.97
1	A	350	VAL	O-C-N	5.09	126.81	121.87
1	B	205	GLU	CB-CA-C	-5.07	104.69	111.89
1	A	356	ILE	O-C-N	-5.06	116.95	121.91
1	B	270	GLY	O-C-N	-5.04	116.14	122.70
1	A	191	VAL	CA-C-N	-5.03	116.04	119.66
1	A	191	VAL	C-N-CA	-5.03	116.04	119.66

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	2834	0	2756	27	0
1	B	2878	0	2815	19	0
2	A	28	0	16	2	0
2	B	14	0	8	0	0
3	A	10	0	0	0	0
3	B	5	0	0	0	0
4	A	292	0	0	6	0
4	B	325	0	0	3	0
All	All	6386	0	5595	48	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 (48) 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:401[B]:2UZ:SAI	2:A:401[B]:2UZ:O	2.20	0.99
1:A:28[B]:MET:HG3	1:A:29:ALA:N	2.04	0.71
1:B:87:ASP:OD2	1:B:92:TYR:OH	2.12	0.68
1:B:262[B]:THR:CG2	1:B:298:VAL:HG12	2.24	0.67
1:A:243:ILE:CD1	1:A:252:ILE:HD12	2.29	0.62
1:A:278[B]:VAL:HG21	1:A:283:ILE:HD11	1.82	0.62
1:A:262:THR:HG22	4:A:730:HOH:O	2.00	0.59
1:A:243:ILE:CD1	1:A:252:ILE:CD1	2.80	0.59
1:A:265[A]:MET:SD	4:A:730:HOH:O	2.58	0.56
1:A:262:THR:O	1:A:265[A]:MET:HG3	2.06	0.55
1:A:175[B]:GLN:HG3	1:A:180:GLN:OE1	2.08	0.54
1:A:33:ILE:HD12	1:A:33:ILE:N	2.23	0.53
1:A:243:ILE:HD12	1:A:252:ILE:HD12	1.91	0.53
1:B:296[B]:ARG:NH1	4:B:646:HOH:O	2.37	0.53
1:A:276:TRP:CE3	1:A:278[A]:VAL:HG12	2.44	0.52
1:A:87:ASP:OD2	1:A:92:TYR:OH	2.21	0.52
1:B:243:ILE:CD1	1:B:252:ILE:HD12	2.40	0.52
1:A:334:LEU:HG	1:A:357:LEU:HD22	1.93	0.51
1:A:262:THR:HG23	1:A:265[A]:MET:HG3	1.94	0.49
1:A:316[A]:THR:HG22	4:A:534:HOH:O	2.14	0.48
1:B:26:PRO:HB3	1:B:48:ILE:HD11	1.95	0.48
1:A:262:THR:CG2	1:A:265[A]:MET:SD	3.02	0.47
1:A:119:LEU:HA	1:A:151:ALA:HA	1.97	0.46
1:B:276:TRP:CD2	1:B:277:PRO:HA	2.51	0.46
1:B:262[B]:THR:HG23	1:B:298:VAL:HG12	1.95	0.46
1:A:344:TYR:CZ	1:A:349:ARG:HG2	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:293:LEU:C	1:A:293:LEU:HD23	2.41	0.45
1:B:334:LEU:HG	1:B:357:LEU:HD22	1.98	0.45
2:A:401[B]:2UZ:H6	4:A:766:HOH:O	2.15	0.44
1:B:243:ILE:CD1	1:B:252:ILE:CD1	2.95	0.44
1:A:276:TRP:CD2	1:A:277:PRO:HA	2.52	0.44
1:A:262:THR:HG23	1:A:265[A]:MET:CG	2.48	0.43
1:A:17:THR:HB	1:A:18:PRO:HD3	2.00	0.43
1:B:7[A]:GLN:HG2	1:B:8:ILE:N	2.33	0.43
1:B:33:ILE:HD12	1:B:33:ILE:N	2.34	0.42
1:A:240:PRO:O	1:A:249:GLN:HG3	2.19	0.42
1:A:204:ARG:NH2	4:A:546:HOH:O	2.52	0.42
1:A:262:THR:HG22	1:A:265[A]:MET:SD	2.60	0.42
1:B:265[A]:MET:HG3	4:B:504:HOH:O	2.20	0.42
1:A:256:GLN:HG3	4:A:639:HOH:O	2.19	0.42
1:B:262[B]:THR:HG22	1:B:298:VAL:HG12	1.99	0.42
1:B:344:TYR:CE1	1:B:349:ARG:HG2	2.55	0.41
1:B:278[A]:VAL:HG22	1:B:354:TRP:CZ3	2.55	0.41
1:A:125:VAL:HG11	1:A:131[A]:LEU:HD23	2.03	0.41
1:B:262[B]:THR:CG2	1:B:298:VAL:CG1	2.96	0.41
1:B:290:LYS:NZ	4:B:700:HOH:O	2.52	0.41
1:B:120[B]:GLN:NE2	1:B:152:ASN:HD22	2.17	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	363/358 (101%)	358 (99%)	5 (1%)	0	100	100
1	B	369/358 (103%)	363 (98%)	6 (2%)	0	100	100
All	All	732/716 (102%)	721 (98%)	11 (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	292/292 (100%)	289 (99%)	3 (1%)	68	43
1	B	299/292 (102%)	295 (99%)	4 (1%)	61	33
All	All	591/584 (101%)	584 (99%)	7 (1%)	65	36

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	293	LEU
1	A	351	ASP
1	A	361	GLN
1	B	205	GLU
1	B	262[A]	THR
1	B	262[B]	THR
1	B	309	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	102	ASN
1	B	346	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

6 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
3	PO4	A	402[B]	-	4,4,4	1.87	1 (25%)	6,6,6	0.50	0
3	PO4	B	402[B]	-	4,4,4	1.56	1 (25%)	6,6,6	1.25	1 (16%)
3	PO4	A	403	-	4,4,4	0.47	0	6,6,6	1.52	1 (16%)
2	2UZ	A	401[A]	-	10,14,14	3.45	6 (60%)	9,19,19	1.73	4 (44%)
2	2UZ	B	401	-	10,14,14	3.59	3 (30%)	9,19,19	2.49	5 (55%)
2	2UZ	A	401[B]	-	10,14,14	2.81	5 (50%)	9,19,19	1.57	1 (11%)

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	2UZ	B	401	-	-	4/8/22/22	0/0/1/1
2	2UZ	A	401[A]	-	-	4/8/22/22	0/0/1/1
2	2UZ	A	401[B]	-	-	6/8/22/22	0/0/1/1

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	2UZ	CAG-CAF	-8.75	1.38	1.49
2	A	401[A]	2UZ	CAG-CAF	-8.72	1.38	1.49
2	A	401[B]	2UZ	CAG-CAF	-6.42	1.41	1.49
2	B	401	2UZ	CA-C	5.63	1.63	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401[A]	2UZ	CAL-CAJ	-3.15	1.41	1.48
2	B	401	2UZ	CAF-CAL	3.08	1.42	1.34
2	A	401[B]	2UZ	CAF-CAL	2.99	1.42	1.34
2	A	401[B]	2UZ	CB-SAI	-2.82	1.75	1.82
2	A	401[A]	2UZ	OAD-CAJ	2.73	1.29	1.22
2	A	401[A]	2UZ	CAF-CAL	2.60	1.41	1.34
2	A	401[A]	2UZ	OXT-C	-2.60	1.22	1.30
2	A	401[B]	2UZ	OAD-CAJ	2.55	1.28	1.22
3	B	402[B]	PO4	P-O3	-2.48	1.47	1.54
3	A	402[B]	PO4	P-O3	-2.36	1.47	1.54
2	A	401[B]	2UZ	CA-C	2.34	1.57	1.52
2	A	401[A]	2UZ	CA-N	2.08	1.53	1.47

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	2UZ	OAB-CAJ-CAL	4.55	124.08	114.27
2	B	401	2UZ	OAB-CAJ-OAD	-3.59	115.36	123.90
2	B	401	2UZ	CAG-SAI-CB	-3.23	88.45	94.35
2	A	401[A]	2UZ	CAF-CAL-NAH	2.54	126.90	120.31
2	A	401[A]	2UZ	C-CA-N	-2.48	104.29	109.49
2	A	401[A]	2UZ	OAB-CAJ-CAL	2.45	119.56	114.27
2	A	401[B]	2UZ	OAB-CAJ-OAD	-2.42	118.14	123.90
2	A	401[A]	2UZ	CAG-SAI-CB	2.41	98.75	94.35
3	A	403	PO4	O3-P-O2	2.39	115.34	107.91
2	B	401	2UZ	OXT-C-O	-2.28	118.90	124.08
3	B	402[B]	PO4	O4-P-O2	2.15	114.61	107.91
2	B	401	2UZ	OXT-C-CA	2.11	121.43	114.15

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401[B]	2UZ	OAB-CAJ-CAL-NAH
2	B	401	2UZ	OAD-CAJ-CAL-NAH
2	B	401	2UZ	OAB-CAJ-CAL-NAH
2	A	401[B]	2UZ	OXT-C-CA-N
2	B	401	2UZ	OAB-CAJ-CAL-CAF
2	A	401[A]	2UZ	O-C-CA-CB
2	A	401[A]	2UZ	OXT-C-CA-N
2	A	401[A]	2UZ	OXT-C-CA-CB
2	A	401[B]	2UZ	OAD-CAJ-CAL-CAF

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Mol	Chain	Res	Type	Atoms
2	A	401[B]	2UZ	OAB-CAJ-CAL-CAF
2	B	401	2UZ	OAD-CAJ-CAL-CAF
2	A	401[B]	2UZ	OAD-CAJ-CAL-NAH
2	A	401[A]	2UZ	O-C-CA-N
2	A	401[B]	2UZ	O-C-CA-N

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401[B]	2UZ	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	351/358 (98%)	-0.12	10 (2%) 55 59	6, 16, 33, 48	16 (4%)
1	B	358/358 (100%)	-0.31	3 (0%) 82 86	6, 15, 28, 48	12 (3%)
All	All	709/716 (99%)	-0.21	13 (1%) 67 72	6, 16, 30, 48	28 (3%)

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	281	ASP	4.2
1	A	279	ASN	3.8
1	A	4	ALA	3.5
1	A	280	PRO	2.9
1	B	241	LEU	2.6
1	B	244	ASN	2.6
1	A	291	ILE	2.5
1	A	126	LYS	2.5
1	A	292	ALA	2.4
1	A	310	ALA	2.4
1	A	283	ILE	2.1
1	B	354	TRP	2.1
1	A	278[A]	VAL	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 oligosaccharides 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. 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	B-factors(Å ²)	Q<0.9
2	2UZ	B	401	14/14	0.81	0.21	24,58,83,91	0
2	2UZ	A	401[B]	14/14	0.85	0.14	19,37,49,49	14
2	2UZ	A	401[A]	14/14	0.85	0.14	23,32,36,41	14
3	PO4	A	403	5/5	0.90	0.14	27,27,33,36	5
3	PO4	B	402[B]	5/5	0.92	0.13	18,19,21,22	5
3	PO4	A	402[B]	5/5	0.97	0.07	14,14,16,18	5

6.5 Other polymers [i](#)

There are no such residues in this entry.